L Number	Hits	Search Text	DB	Time stamp
1	322	(548/468).CCLS.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/08/24 14:28
2	410	(548/486).CCLS.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/08/24 14:28
3	1029	(514/415).CCLS.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/08/24 14:29
4	190	tang.inv. adj peng.inv.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/08/24 14:29

L Number	Hits	Search Text	DB	Time stamp
1	322	(548/468).CCLS.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/08/24 14:28
2	410	(548/486).CCLS.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/08/24 14:28
3	1029	(514/415).CCLS.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/08/24 14:29
4	190	tang.inv. adj peng.inv.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/08/24 14:29

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NEWS 12	AUG 02	CAplus and CA patent records enhanced with European and Japan Patent Office Classifications
NEWS 13	AUG 02	STN User Update to be held August 22 in conjunction with the 228th ACS National Meeting
NEWS 14	AUG 02	The Analysis Edition of STN Express with Discover! (Version 7.01 for Windows) now available
NEWS 15	AUG 04	Pricing for the Save Answers for SciFinder Wizard within STN Express with Discover! will change September 1, 2004

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FILE 'HOME' ENTERED AT 13:39:11 ON 24 AUG 2004

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'REGISTRY' ENTERED AT 13:39:23 ON 24 AUG 2004

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 23 AUG 2004 HIGHEST RN 731771-88-3
DICTIONARY FILE UPDATES: 23 AUG 2004 HIGHEST RN 731771-88-3

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:
http://www.cas.org/ONLINE/DBSS/registryss.html

=> ....Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

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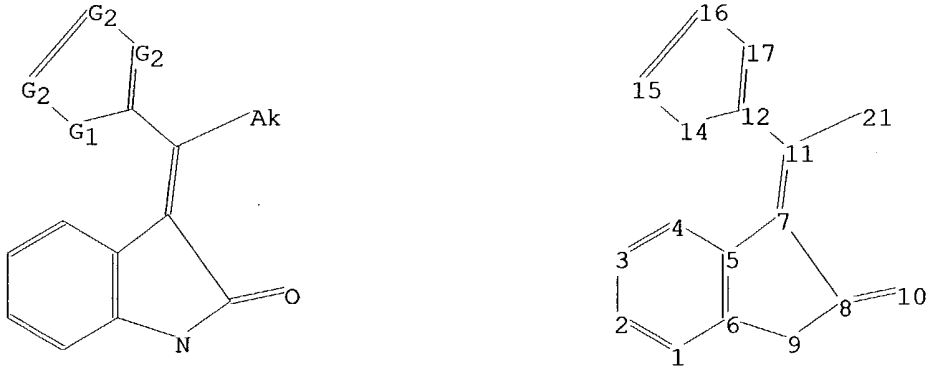
Uploading C:\STNEXP4\QUERIES\10725079-6.str



chain nodes :
10 11 21
ring nodes :
1 2 3 4 5 6 7 8 9 12 14 15 16 17
chain bonds :
7-11 8-10 11-12 11-21

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ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 12-14 12-17 14-15 15-16 16-17

exact/norm bonds :
5-7 6-9 7-8 7-11 8-9 8-10 11-12 11-21 12-14 12-17 14-15 15-16 16-17
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6
isolated ring systems :
containing 1 : 12 :

G1:O,S,N

G2:C,O,S,N

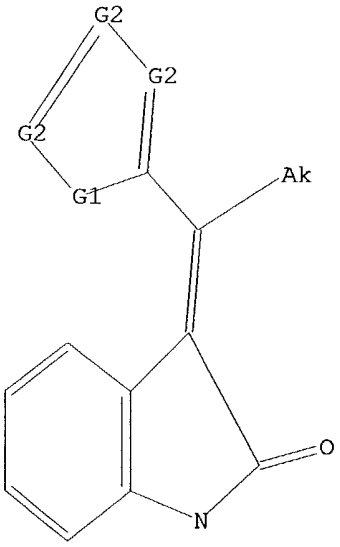
Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS
11:CLASS 12:Atom 14:Atom 15:Atom 16:Atom 17:Atom 21:CLASS

L1 STRUCTURE UPLOADED

=> que L1

L2 QUE L1

=> d l1
L1 HAS NO ANSWERS
L1 STR



G1 O,S,N
G2 C,O,S,N

Structure attributes must be viewed using STN Express query preparation.

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=> s l1 sss sam
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SAMPLE SCREEN SEARCH COMPLETED -     634 TO ITERATE

100.0% PROCESSED      634 ITERATIONS                                 1 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:        11170 TO    14190
PROJECTED ANSWERS:               1 TO       80

L3             1 SEA SSS SAM L1

=> d scan
```

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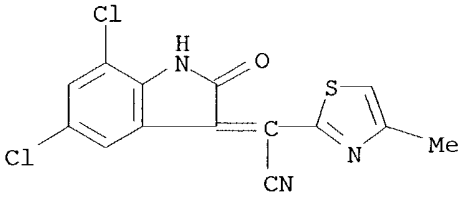
L3 1 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2-Thiazoleacetonitrile, α-(5,7-dichloro-1,2-dihydro-2-oxo-3H-indol-3-ylidene)-4-methyl- (9CI)
MF C14 H7 Cl2 N3 O S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

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=> s l1 sss sam
SAMPLE SEARCH INITIATED 13:40:09 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -     634 TO ITERATE

100.0% PROCESSED     634 ITERATIONS                                 1 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:        11170 TO    14190
PROJECTED ANSWERS:               1 TO       80

L4            1 SEA SSS SAM L1

=> s l1 sss ful
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FULL SCREEN SEARCH COMPLETED -  12602 TO ITERATE

100.0% PROCESSED   12602 ITERATIONS                               23 ANSWERS
SEARCH TIME: 00.00.01

L5           23 SEA SSS FUL L1

=> file caplus
COST IN U.S. DOLLARS                                SINCE FILE      TOTAL
                                                       ENTRY    SESSION
FULL ESTIMATED COST                                   155.84     156.05

FILE 'CAPLUS' ENTERED AT 13:40:22 ON 24 AUG 2004
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FILE LAST UPDATED: 23 Aug 2004  (20040823/ED)

  This file contains CAS Registry Numbers for easy and accurate
  substance identification.


=> s l5
L6            3 L5

=> d l6 1-3 bib hitstr
```
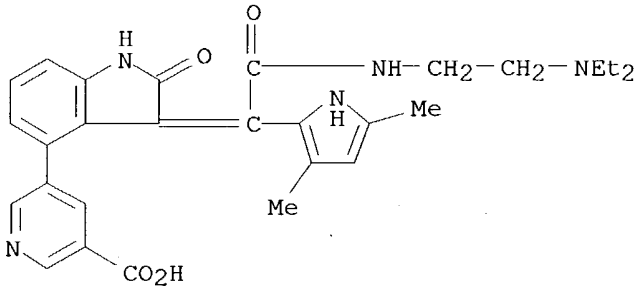

L6 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN
AN 2002:31440 CAPLUS
DN 136:102386
TI Preparation and use of 4-heteroaryl-3-heteroarylidenyl-2-indolinones and their use as protein kinase inhibitors
IN Tang, Peng Cho; Wei, Chung Chen; Huang, Ping; Cui, Jingron
PA Sugen, Inc., USA
SO PCT Int. Appl., 164 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2002002551	A1	20020110	WO 2001-US20768	20010629
	W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
	US 2002187978	A1	20021212	US 2001-894902	20010629
	US 6635640	B2	20031021		
	EP 1296975	A1	20030402	EP 2001-948830	20010629
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR				
	JP 2004502686	T2	20040129	JP 2002-507803	20010629
	US 2004097497	A1	20040520	US 2003-648810	20030827
PRAI	US 2000-215654P	P	20000630		
	US 2001-894902	A3	20010629		
	WO 2001-US20768	W	20010629		

OS MARPAT 136:102386
IT **388117-27-9P**, 5-[3-[4-(2-Diethylaminoethylcarbamoyl)-3,5-dimethyl-1H-pyrrol-2-ylmethylene]-2-oxo-2,3-dihydro-1H-indol-4-yl]nicotinic acid
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(drug; preparation and use of 4-heteroaryl-3-heteroarylidenyl-2-indolinones and their use as protein kinase inhibitors)
RN 388117-27-9 CAPLUS
CN 3-Pyridinecarboxylic acid, 5-[3-[2-[[2-(diethylamino)ethyl]amino]-1-(3,5-dimethyl-1H-pyrrol-2-yl)-2-oxoethylidene]-2,3-dihydro-2-oxo-1H-indol-4-yl]- (9CI) (CA INDEX NAME)



RE.CNT 9 THERE ARE 9 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

V. Balasubramanian

L6 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN
AN 2001:816637 CAPLUS
DN 135:344374
TI Preparation of oxindolylidenylacetic acid derivatives and their use as protein kinase inhibitors
IN Wei, Chung-Chen
PA Sugen, Inc., USA
SO PCT Int. Appl., 99 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2001083450	A2	20011108	WO 2001-US14230	20010502
	WO 2001083450	A3	20020411		
	W: AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CO, CU, CZ, DE, DK, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
	US 2002028828	A1	20020307	US 2001-846340	20010502
	EP 1299355	A2	20030409	EP 2001-929000	20010502
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR				
	JP 2003531895	T2	20031028	JP 2001-580879	20010502
	US 2003216462	A1	20031120	US 2003-371157	20030224
	US 2004039196	A1	20040226	US 2003-460641	20030613
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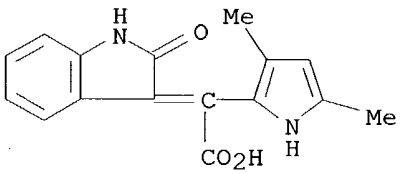
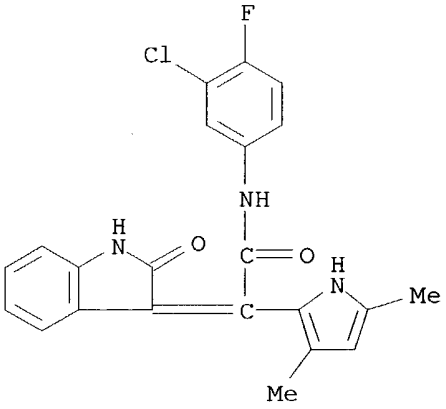
	US 2001-846340	B1	20010502		
	WO 2001-US14230	W	20010502		
	US 2003-371157	A3	20030224		

OS MARPAT 135:344374
IT **371786-23-1P 371786-25-3P 371786-26-4P**
371786-27-5P 371786-28-6P 371786-29-7P
371786-30-0P 371786-31-1P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(drug; preparation of oxindolylidenylacetic acid derivs. and their use as protein kinase inhibitors)
RN 371786-23-1 CAPLUS
CN 1H-Pyrrole-2-acetic acid, α-(1,2-dihydro-2-oxo-3H-indol-3-ylidene)-3,5-dimethyl- (9CI) (CA INDEX NAME)



RN 371786-25-3 CAPLUS
CN 1H-Pyrrole-2-acetamide, N-(3-chloro-4-fluorophenyl)-α-(1,2-dihydro-2-oxo-3H-indol-3-ylidene)-3,5-dimethyl- (9CI) (CA INDEX NAME)



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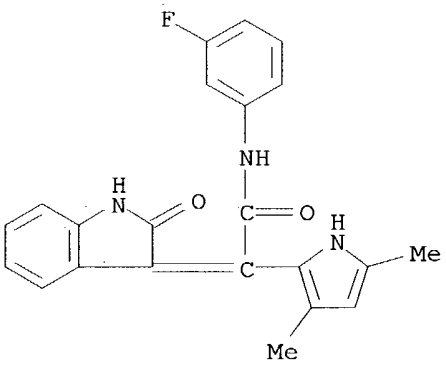
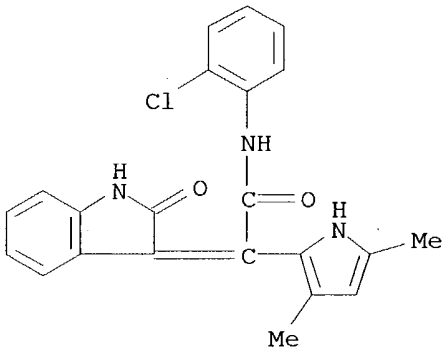
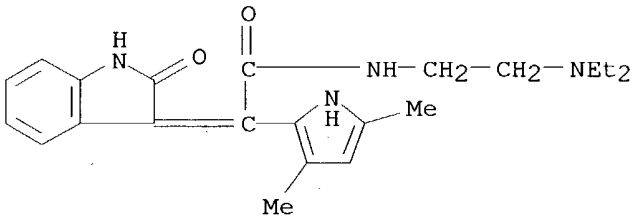


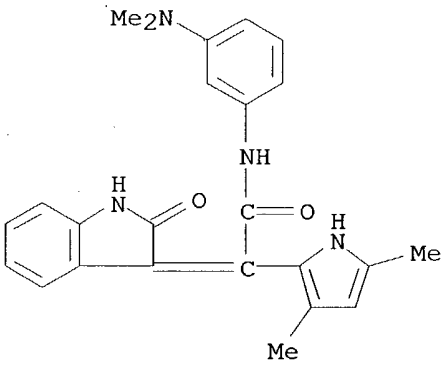
RN 371786-27-5 CAPLUS
CN 1H-Pyrrole-2-acetamide, N-(2-chlorophenyl)-α-(1,2-dihydro-2-oxo-3H-indol-3-ylidene)-3,5-dimethyl- (9CI) (CA INDEX NAME)



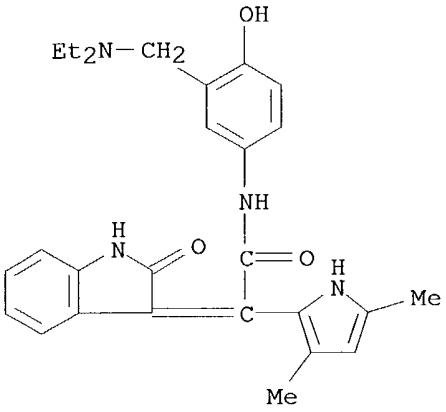
RN 371786-28-6 CAPLUS
CN 1H-Pyrrole-2-acetamide, N-[2-(diethylamino)ethyl]-α-(1,2-dihydro-2-oxo-3H-indol-3-ylidene)-3,5-dimethyl- (9CI) (CA INDEX NAME)



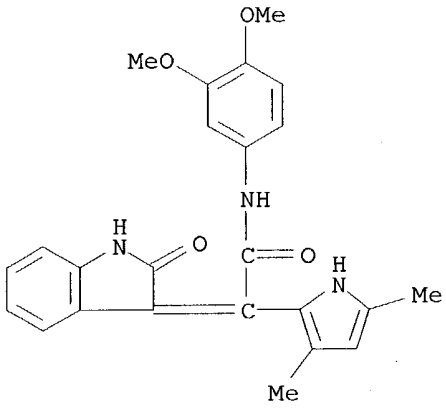
RN 371786-29-7 CAPLUS
CN 1H-Pyrrole-2-acetamide, α-(1,2-dihydro-2-oxo-3H-indol-3-ylidene)-N-[3-(dimethylamino)phenyl]-3,5-dimethyl- (9CI) (CA INDEX NAME)



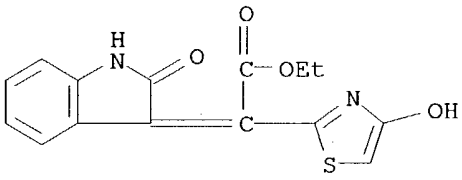
RN 371786-30-0 CAPLUS
CN 1H-Pyrrole-2-acetamide, N-[3-[(diethylamino)methyl]-4-hydroxyphenyl]-α-(1,2-dihydro-2-oxo-3H-indol-3-ylidene)-3,5-dimethyl- (9CI) (CA INDEX NAME)



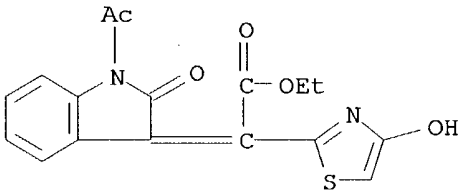
RN 371786-31-1 CAPLUS
CN 1H-Pyrrole-2-acetamide, α-(1,2-dihydro-2-oxo-3H-indol-3-ylidene)-N-(3,4-dimethoxyphenyl)-3,5-dimethyl- (9CI) (CA INDEX NAME)



L6 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2004 ACS on STN
AN 1987:617539 CAPLUS
DN 107:217539
TI Nitriles in heterocyclic synthesis. A novel synthesis of spiropyran-4-ylindolidene derivatives
AU Hafez, Ebtisam Abdel Aziz; Abdul Galil, Fathy M.; Sherif, Sherif M.; Elnagdi, Mohamed H.
CS Fac. Sci., Cairo Univ., Giza, Egypt
SO Journal of Heterocyclic Chemistry (1986), 23(5), 1375-8
CODEN: JHTCAD; ISSN: 0022-152X
DT Journal
LA English
OS CASREACT 107:217539
IT **111277-27-1P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation and reactions of)
RN 111277-27-1 CAPLUS
CN 2-Thiazoleacetic acid, α-(1,2-dihydro-2-oxo-3H-indol-3-ylidene)-4-hydroxy-, ethyl ester (9CI) (CA INDEX NAME)



IT **111348-06-2P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)
RN 111348-06-2 CAPLUS
CN 2-Thiazoleacetic acid, α-(1-acetyl-1,2-dihydro-2-oxo-3H-indol-3-ylidene)-4-hydroxy-, ethyl ester (9CI) (CA INDEX NAME)



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=> log y

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	10.25	166.30

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NEWS 15	AUG 04	Pricing for the Save Answers for SciFinder Wizard within STN Express with Discover! will change September 1, 2004

NEWS EXPRESS	JULY 30 CURRENT WINDOWS VERSION IS V7.01, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 11 AUGUST 2004
NEWS HOURS	STN Operating Hours Plus Help Desk Availability
NEWS INTER	General Internet Information
NEWS LOGIN	Welcome Banner and News Items
NEWS PHONE	Direct Dial and Telecommunication Network Access to STN
NEWS WWW	CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

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* * * * * * * * * * * * * STN Columbus * * * * * * * * * * * * * * *

V. Balasubramanian

FILE 'HOME' ENTERED AT 13:22:38 ON 24 AUG 2004

=> file reg
COST IN U.S. DOLLARS

	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'REGISTRY' ENTERED AT 13:22:52 ON 24 AUG 2004
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

Property values tagged with IC are from the ZIC/VINITI data file
provided by InfoChem.

STRUCTURE FILE UPDATES: 23 AUG 2004 HIGHEST RN 731771-88-3
DICTIONARY FILE UPDATES: 23 AUG 2004 HIGHEST RN 731771-88-3

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

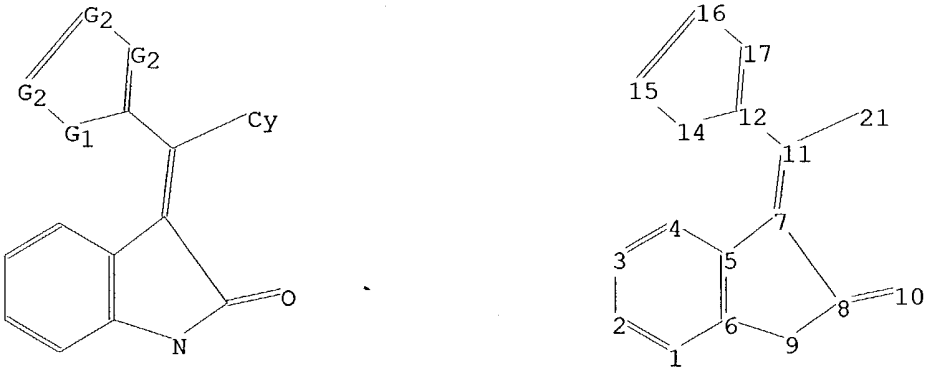
Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
information enter HELP PROP at an arrow prompt in the file or refer
to the file summary sheet on the web at:
http://www.cas.org/ONLINE/DBSS/registryss.html

=> ....Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=>
Uploading C:\STNEXP4\QUERIES\10725079-5.str



chain nodes :
10 11 21
ring nodes :
1 2 3 4 5 6 7 8 9 12 14 15 16 17
chain bonds :
7-11 8-10 11-12 11-21

V. Balasubramanian

ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 12-14 12-17 14-15 15-16 16-17

exact/norm bonds :
5-7 6-9 7-8 7-11 8-9 8-10 11-12 11-21 12-14 12-17 14-15 15-16 16-17
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6
isolated ring systems :
containing 1 : 12 :

G1:O,S,N

G2:C,O,S,N

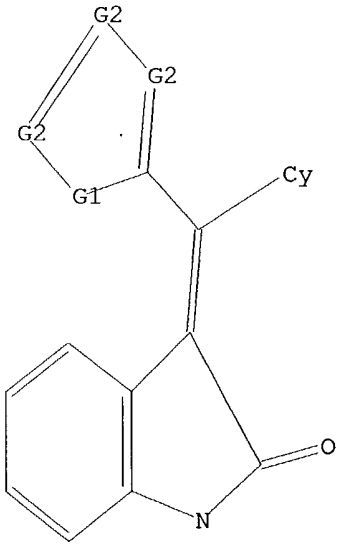
Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS
11:CLASS 12:Atom 14:Atom 15:Atom 16:Atom 17:Atom 21:CLASS

L1 STRUCTURE UPLOADED

=> que L1

L2 QUE L1

=> d l1
L1 HAS NO ANSWERS
L1 STR



G1 O,S,N
G2 C,O,S,N

Structure attributes must be viewed using STN Express query preparation.

V. Balasubramanian

```
=> s l1 sss sam
SAMPLE SEARCH INITIATED 13:23:12 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -     634 TO ITERATE

100.0% PROCESSED      634 ITERATIONS                                  2 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:         11170 TO    14190
PROJECTED ANSWERS:                2 TO      124

L3             2 SEA SSS SAM L1

=> d scan
```

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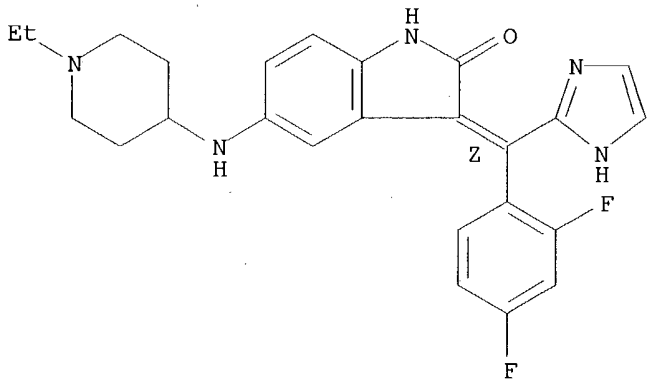
L3 2 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 3-[(2,4-difluorophenyl)-1H-imidazol-2-ylmethylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI)
MF C25 H25 F2 N5 O

Double bond geometry as shown.



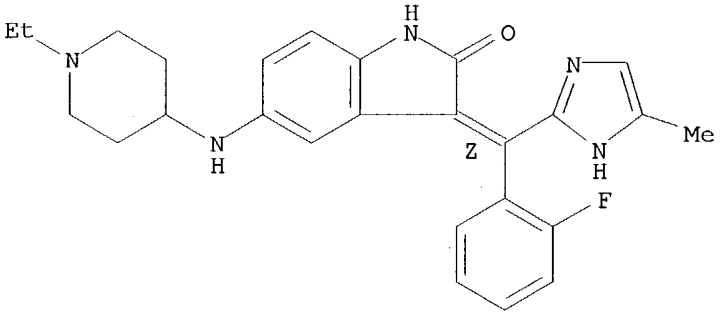
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

V. Balasubramanian

L3 2 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[(2-fluorophenyl)(4-methyl-1H-imidazol-2-yl)methylene]-1,3-dihydro-, (3Z)- (9CI)
MF C26 H28 F N5 O

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

V. Balasubramanian

=> s l1 sss ful
FULL SEARCH INITIATED 13:23:37 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 12602 TO ITERATE

100.0% PROCESSED 12602 ITERATIONS 73 ANSWERS
SEARCH TIME: 00.00.01

L4 73 SEA SSS FUL L1

=> file caplus
COST IN U.S. DOLLARS SINCE FILE ENTRY TOTAL SESSION
FULL ESTIMATED COST 155.42 155.63

FILE 'CAPLUS' ENTERED AT 13:23:43 ON 24 AUG 2004
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

FILE COVERS 1907 - 24 Aug 2004 VOL 141 ISS 9
FILE LAST UPDATED: 23 Aug 2004 (20040823/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l4
L5 2 L4

=> d l5 1-2 bib hitstr

L5 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2004 ACS on STN
AN 2004:493723 CAPLUS
DN 141:54195
TI Preparation of oxindole derivatives as kinase modulators
IN Bannen, Lynne Canne; Brown, S. David; Cheng, Wei; Co, Erick Wang; Nuss, John M.; Kim, Moon Hwan; Klein, Rhett Ronald; Le, Donna T.; Lew, Amy; Mac, Morrison B.; Parks, Jason Jevious; Wen, Zhaoyang; Xu, Wei
PA Exelixis, Inc., USA
SO PCT Int. Appl., 120 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2004050681	A2	20040617	WO 2003-US36567	20031114

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW, AM, AZ, BY, KG, KZ, MD
RW: BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG

PRAI	US 2002-426680P	P	20021115
	US 2003-470674P	P	20030514

OS MARPAT 141:54195
IT **705945-81-9P 705945-90-0P 705945-92-2P**
705946-05-0P 705946-16-3P 705946-17-4P
705946-20-9P 705946-24-3P 705946-25-4P
705946-27-6P 705946-29-8P 705946-30-1P
705946-33-4P 705946-34-5P 705946-35-6P
705946-36-7P 705946-39-0P 705946-40-3P
705946-41-4P 705946-42-5P 705946-43-6P
705946-44-7P 705946-45-8P 705946-46-9P
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705946-69-6P 705946-70-9P 705946-71-0P
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705946-75-4P 705946-76-5P 705946-77-6P
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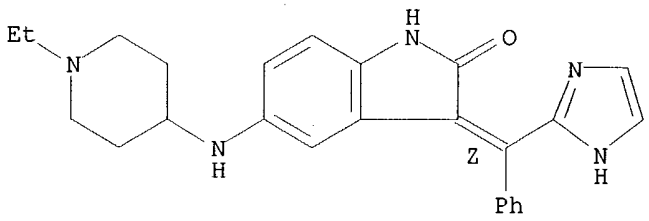
705946-87-8P 705946-88-9P 705946-89-0P
705946-90-3P 705946-91-4P 705946-92-5P
705946-93-6P 705946-94-7P
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of oxindole derivs. as kinase modulators)

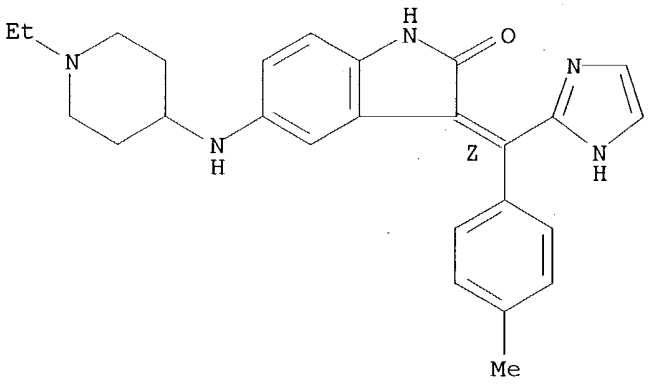
RN 705945-81-9 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-3-(1H-imidazol-2-ylphenylmethylene)-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



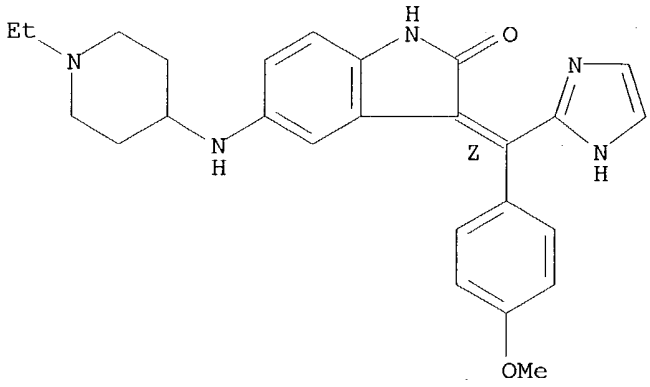
RN 705945-90-0 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-3-[1H-imidazol-2-yl(4-methylphenyl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



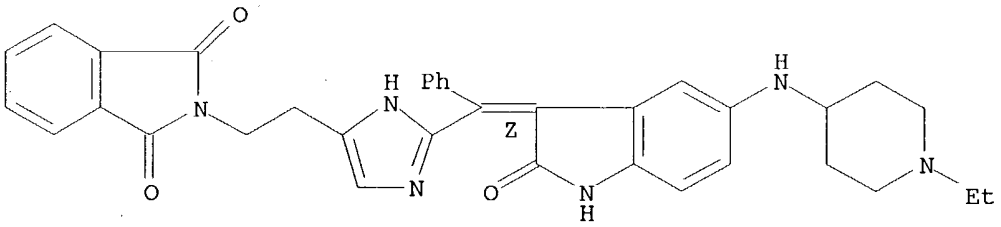
RN 705945-92-2 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-3-[1H-imidazol-2-yl(4-methoxyphenyl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



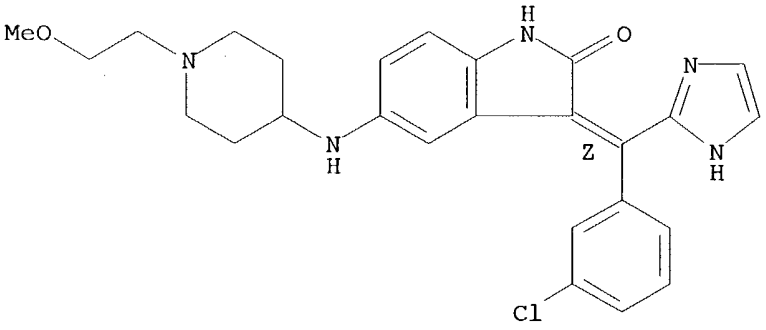
RN 705946-05-0 CAPLUS
CN 1H-Isoindole-1,3(2H)-dione, 2-[2-[2-[(Z)-[5-[(1-ethyl-4-piperidinyl)amino]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]phenylmethyl]-1H-imidazol-4-yl]ethyl]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



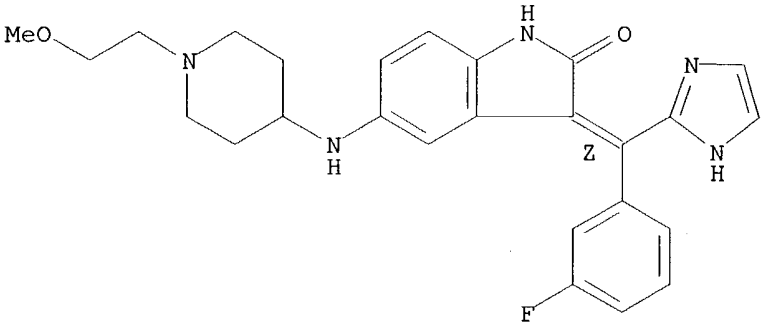
RN 705946-16-3 CAPLUS
CN 2H-Indol-2-one, 3-[(3-chlorophenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



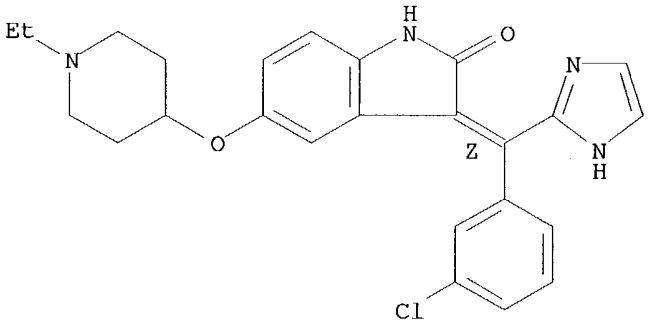
RN 705946-17-4 CAPLUS
CN 2H-Indol-2-one, 3-[(3-fluorophenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



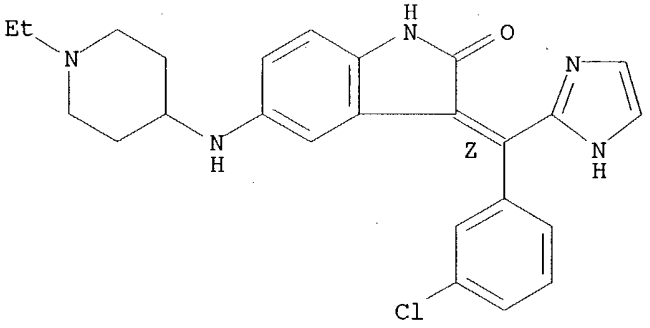
RN 705946-20-9 CAPLUS
CN 2H-Indol-2-one, 3-[(3-chlorophenyl)-1H-imidazol-2-ylmethylene]-5-[(1-ethyl-4-piperidinyl)oxy]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 705946-24-3 CAPLUS
CN 2H-Indol-2-one, 3-[(3-chlorophenyl)-1H-imidazol-2-ylmethylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

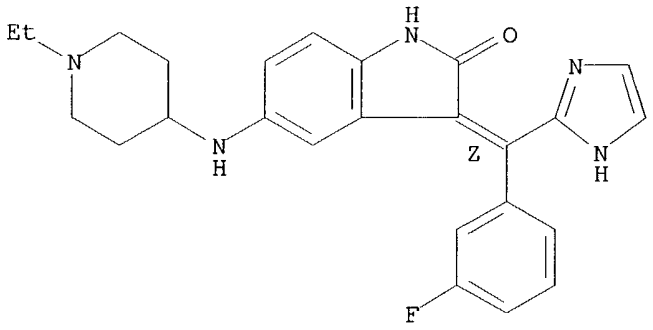
Double bond geometry as shown.



RN 705946-25-4 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[(3-fluorophenyl)-1H-

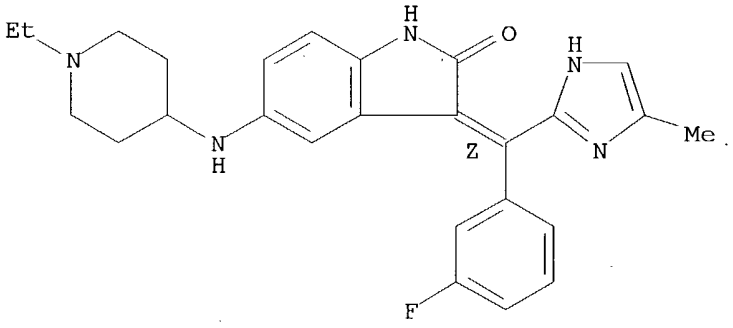
imidazol-2-ylmethylene]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



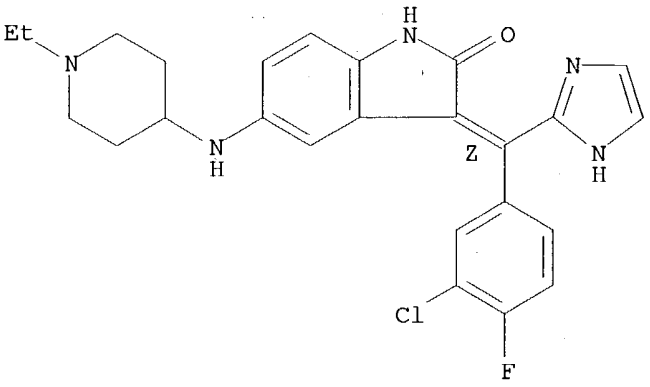
RN 705946-27-6 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[(3-fluorophenyl)(4-methyl-1H-imidazol-2-yl)methylene]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



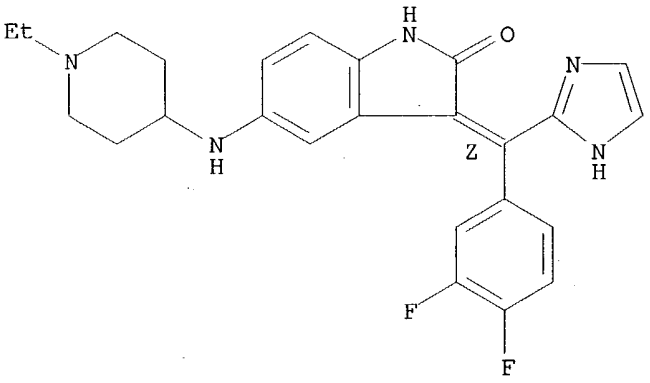
RN 705946-29-8 CAPLUS
CN 2H-Indol-2-one, 3-[(3-chloro-4-fluorophenyl)-1H-imidazol-2-ylmethylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



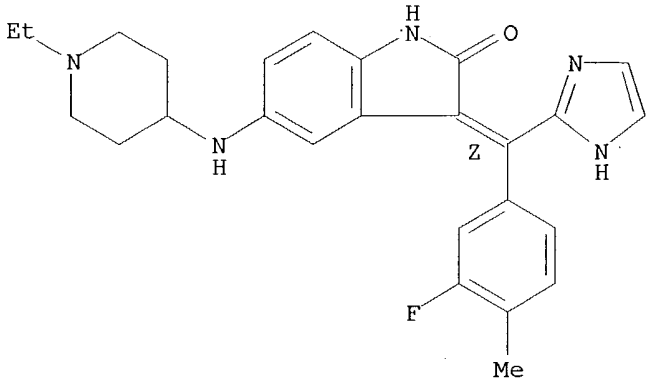
RN 705946-30-1 CAPLUS
CN 2H-Indol-2-one, 3-[(3,4-difluorophenyl)-1H-imidazol-2-ylmethylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



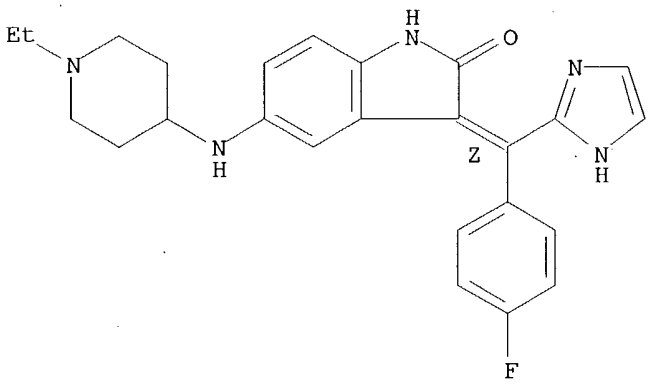
RN 705946-33-4 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[(3-fluoro-4-methylphenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



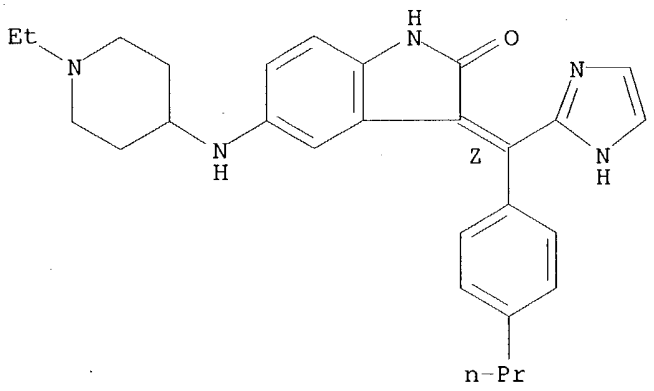
RN 705946-34-5 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[(4-fluorophenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



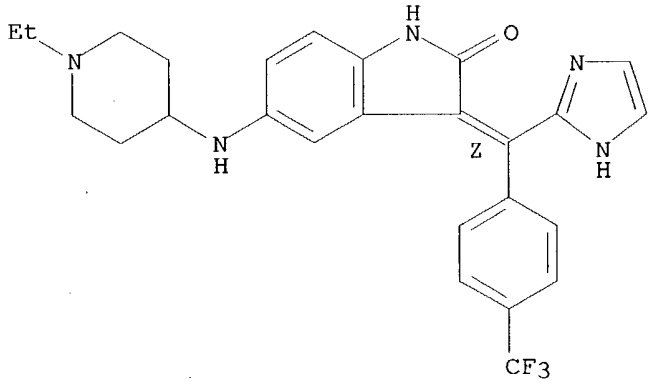
RN 705946-35-6 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-3-[1H-imidazol-2-yl(4-propylphenyl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



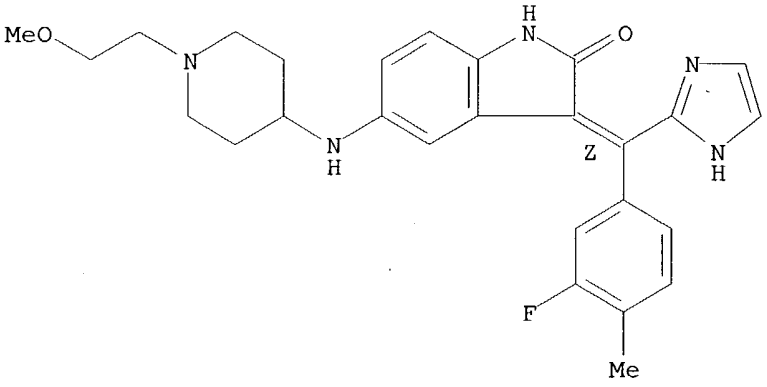
RN 705946-36-7 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-3-[1H-imidazol-2-yl[4-(trifluoromethyl)phenyl]methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



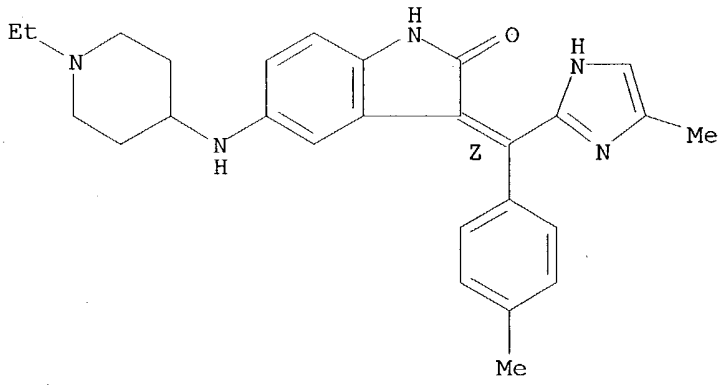
RN 705946-39-0 CAPLUS
CN 2H-Indol-2-one, 3-[(3-fluoro-4-methylphenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



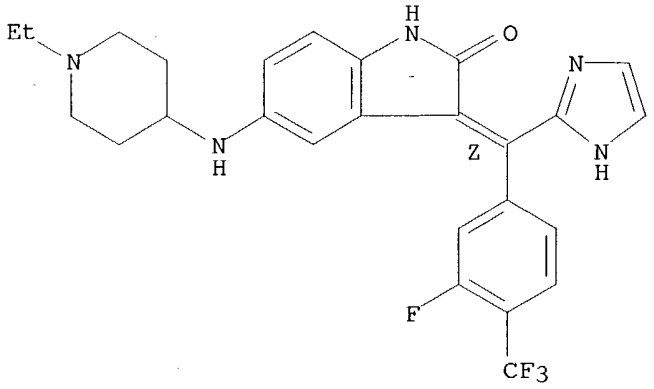
RN 705946-40-3 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-3-[(4-methyl-1H-imidazol-2-yl)(4-methylphenyl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



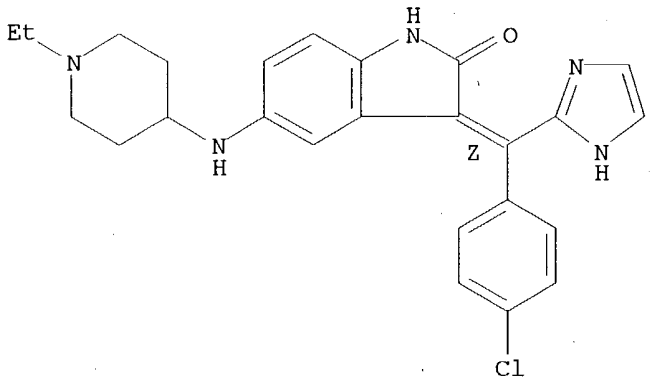
RN 705946-41-4 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[[3-fluoro-4-(trifluoromethyl)phenyl]-1H-imidazol-2-ylmethylene]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



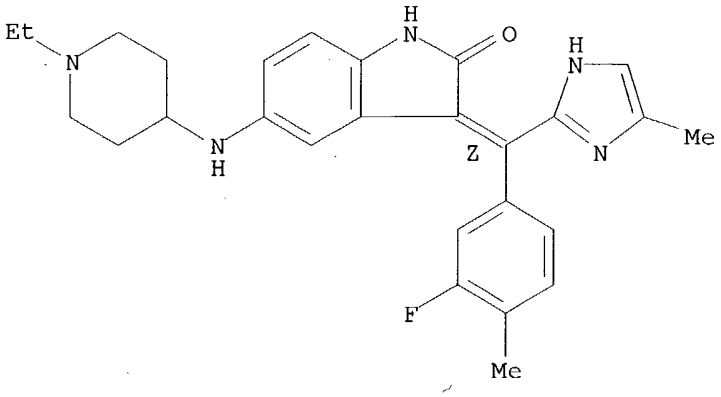
RN 705946-42-5 CAPLUS
CN 2H-Indol-2-one, 3-[(4-chlorophenyl)-1H-imidazol-2-ylmethylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



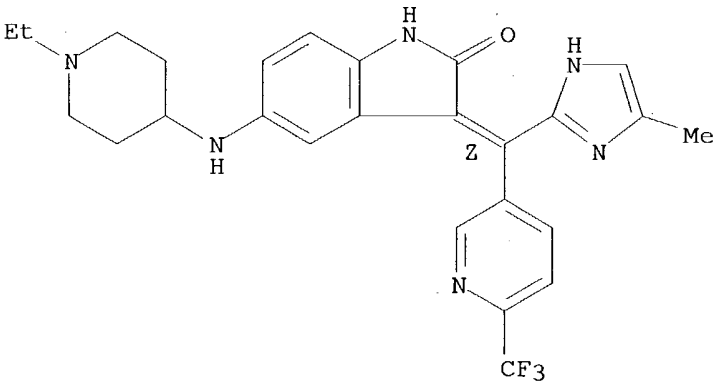
RN 705946-43-6 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[(3-fluoro-4-methylphenyl)(4-methyl-1H-imidazol-2-yl)methylene]-1,3-dihydro-, (3Z)-(9CI) (CA INDEX NAME)

Double bond geometry as shown.



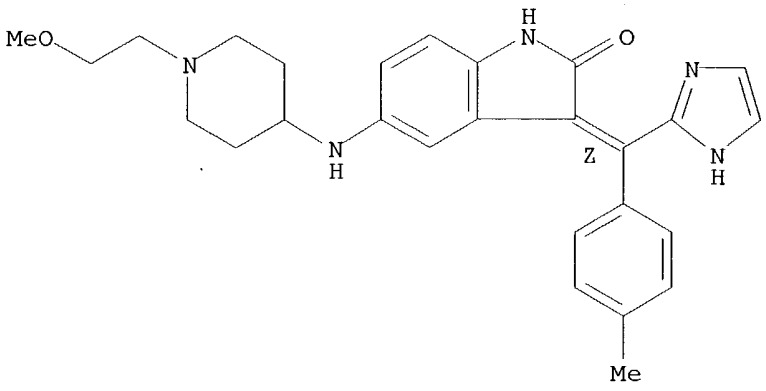
RN 705946-44-7 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-3-[(4-methyl-1H-imidazol-2-yl)[6-(trifluoromethyl)-3-pyridinyl]methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



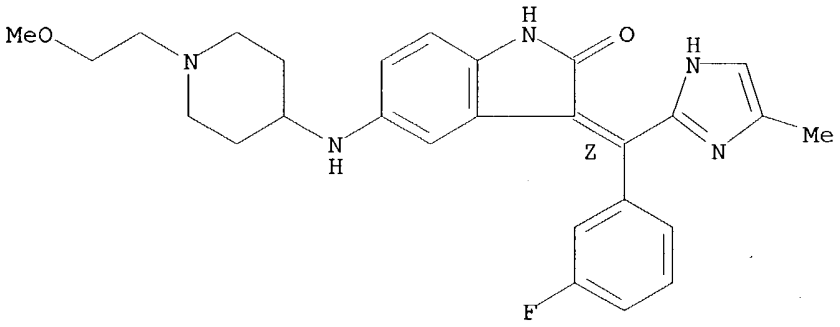
RN 705946-45-8 CAPLUS
CN 2H-Indol-2-one, 1,3-dihydro-3-[1H-imidazol-2-yl(4-methylphenyl)methylene]-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



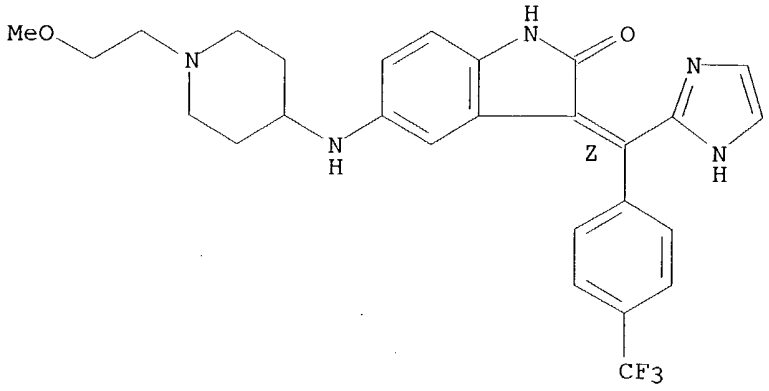
RN 705946-46-9 CAPLUS
CN 2H-Indol-2-one, 3-[(3-fluorophenyl)(4-methyl-1H-imidazol-2-yl)methylene]-1,3-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



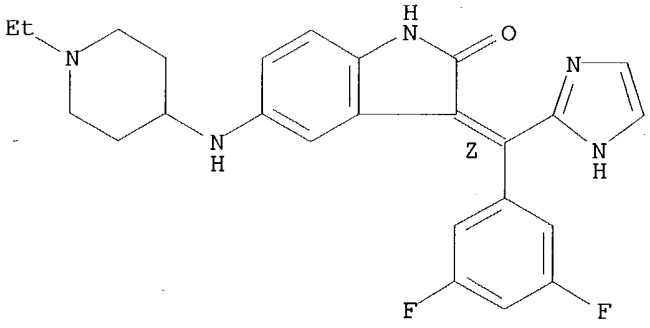
RN 705946-47-0 CAPLUS
CN 2H-Indol-2-one, 1,3-dihydro-3-[1H-imidazol-2-yl[4-(trifluoromethyl)phenyl]methylene]-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



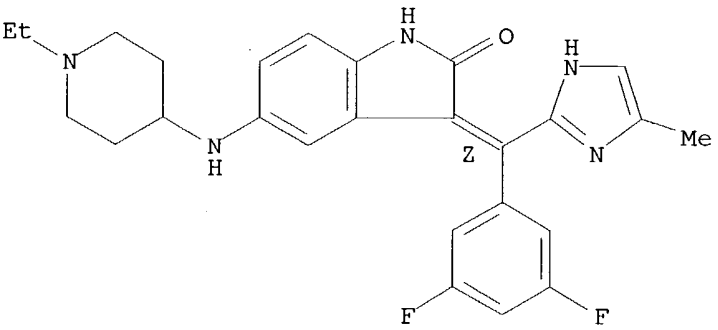
RN 705946-49-2 CAPLUS
CN 2H-Indol-2-one, 3-[(3,5-difluorophenyl)-1H-imidazol-2-ylmethylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



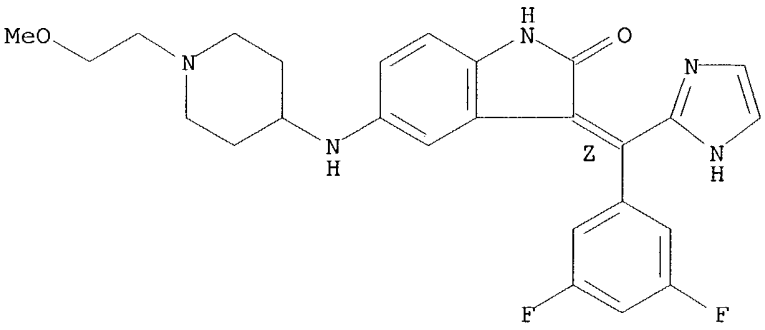
RN 705946-50-5 CAPLUS
CN 2H-Indol-2-one, 3-[(3,5-difluorophenyl)(4-methyl-1H-imidazol-2-yl)methylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



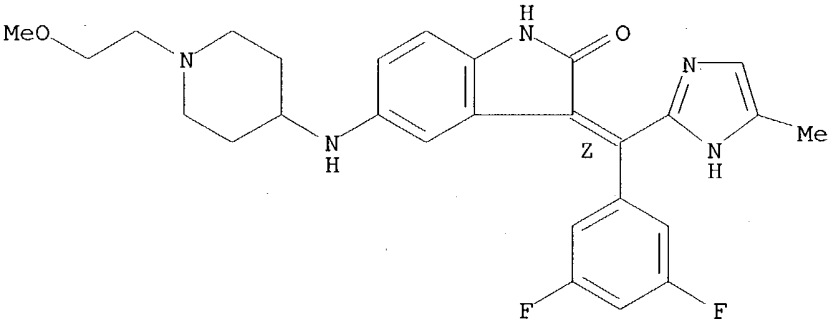
RN 705946-51-6 CAPLUS
CN 2H-Indol-2-one, 3-[(3,5-difluorophenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



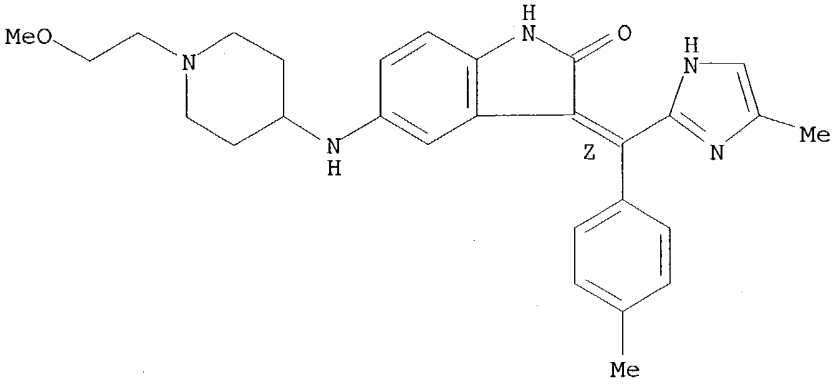
RN 705946-52-7 CAPLUS
CN 2H-Indol-2-one, 3-[(3,5-difluorophenyl)(4-methyl-1H-imidazol-2-yl)methylene]-1,3-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



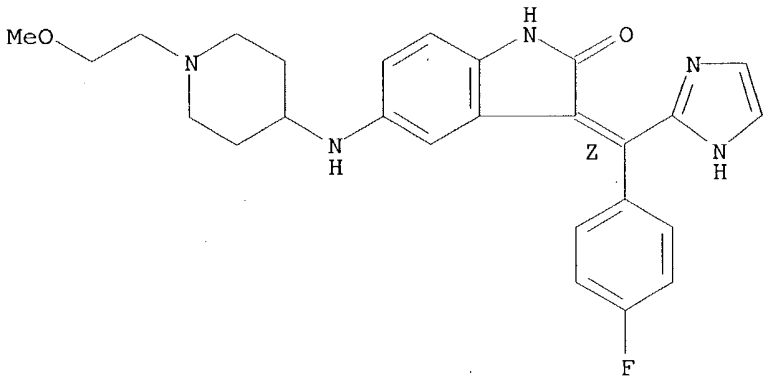
RN 705946-53-8 CAPLUS
CN 2H-Indol-2-one, 1,3-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-3-[(4-methyl-1H-imidazol-2-yl)(4-methylphenyl)methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



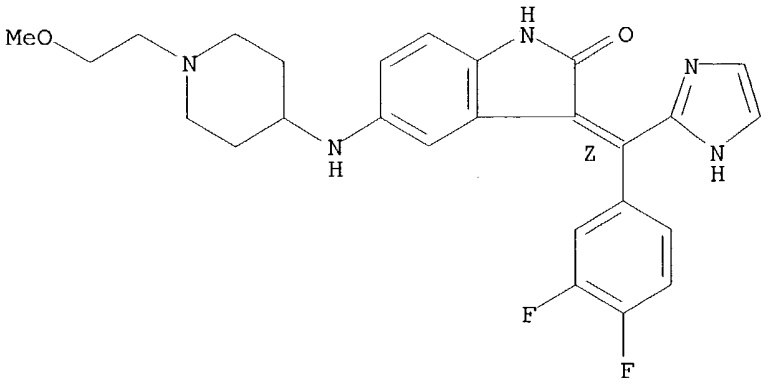
RN 705946-54-9 CAPLUS
CN 2H-Indol-2-one, 3-[(4-fluorophenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



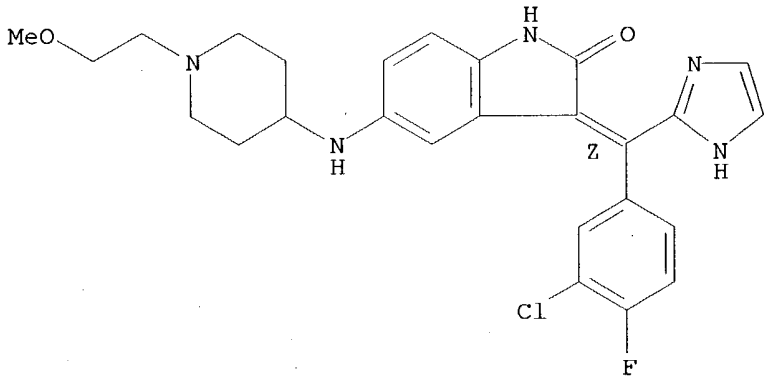
RN 705946-55-0 CAPLUS
CN 2H-Indol-2-one, 3-[(3,4-difluorophenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



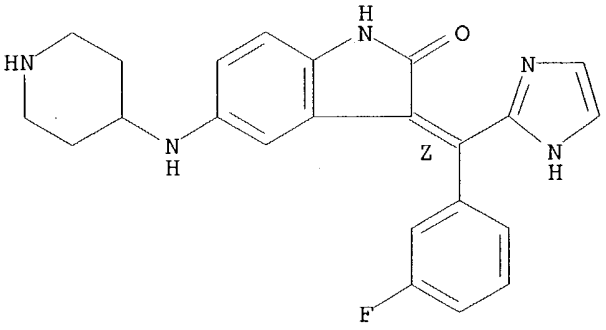
RN 705946-56-1 CAPLUS
CN 2H-Indol-2-one, 3-[(3-chloro-4-fluorophenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



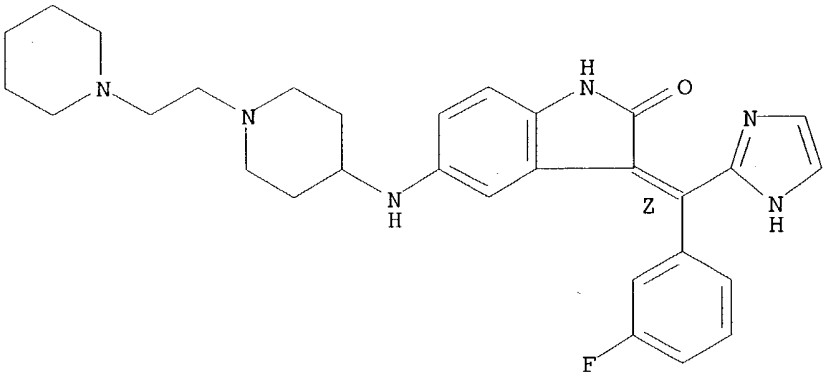
RN 705946-57-2 CAPLUS
CN 2H-Indol-2-one, 3-[(3-fluorophenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-5-(4-piperidinylamino)-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



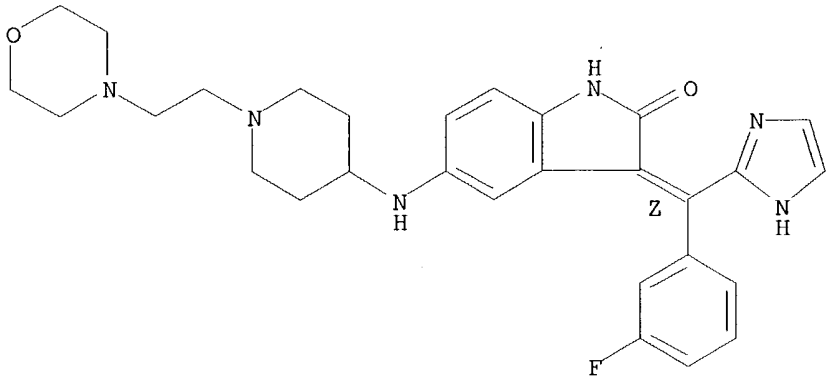
RN 705946-58-3 CAPLUS
CN 2H-Indol-2-one, 3-[(3-fluorophenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-5-[[1-[2-(1-piperidinyl)ethyl]-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



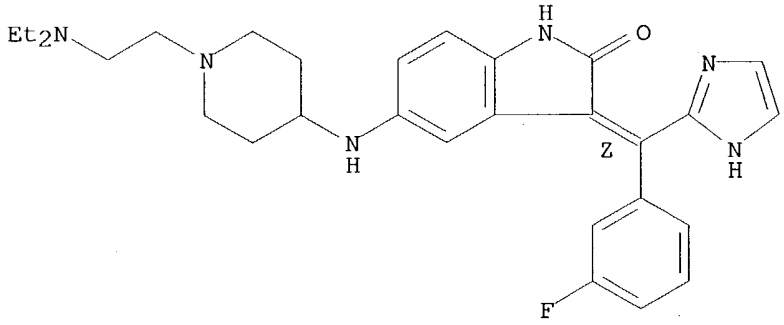
RN 705946-59-4 CAPLUS
CN 2H-Indol-2-one, 3-[(3-fluorophenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-5-[[1-[2-(4-morpholinyl)ethyl]-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



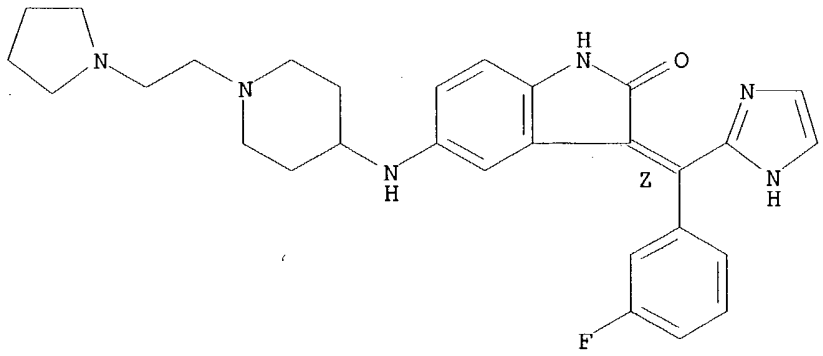
RN 705946-60-7 CAPLUS
CN 2H-Indol-2-one, 5-[[1-[2-(diethylamino)ethyl]-4-piperidinyl]amino]-3-[(3-fluorophenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



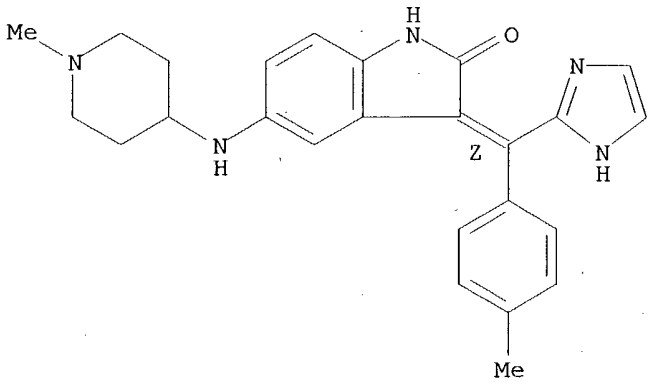
RN 705946-61-8 CAPLUS
CN 2H-Indol-2-one, 3-[(3-fluorophenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-5-[[1-[2-(1-pyrrolidinyl)ethyl]-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



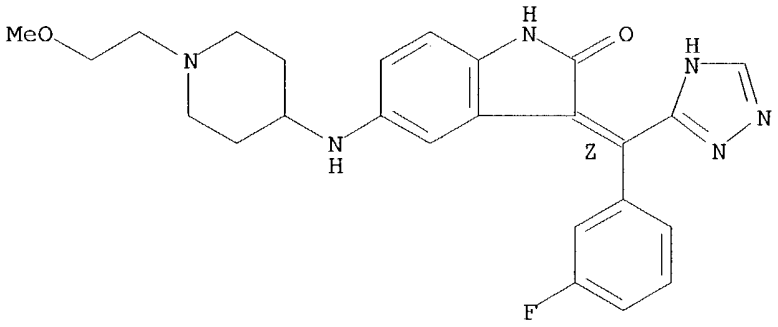
RN 705946-62-9 CAPLUS
CN 2H-Indol-2-one, 1,3-dihydro-3-[1H-imidazol-2-yl(4-methylphenyl)methylene]-5-[(1-methyl-4-piperidinyl)amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



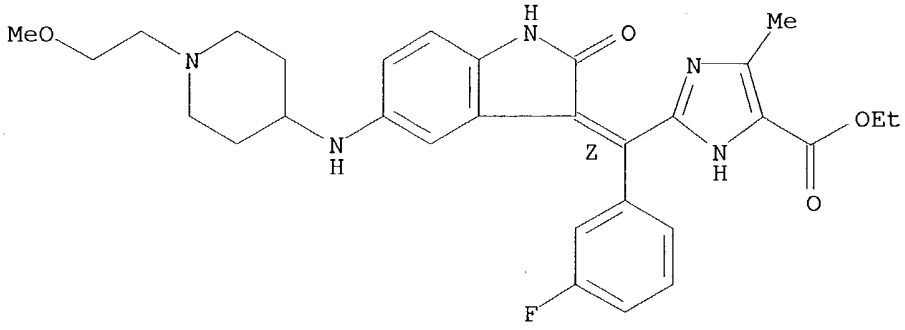
RN 705946-63-0 CAPLUS
CN 2H-Indol-2-one, 3-[(3-fluorophenyl)-1H-1,2,4-triazol-3-ylmethylene]-1,3-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



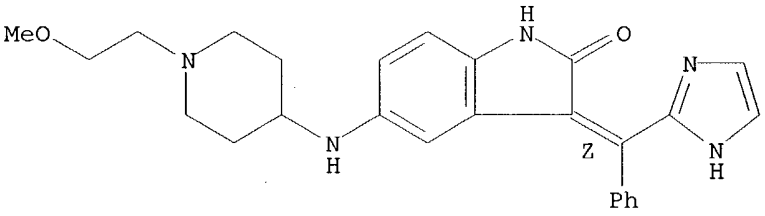
RN 705946-64-1 CAPLUS
CN 1H-Imidazole-4-carboxylic acid, 2-[(Z)-[1,2-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-2-oxo-3H-indol-3-ylidene](3-fluorophenyl)methyl]-5-methyl-, ethyl ester (9CI) (CA INDEX NAME)

Double bond geometry as shown.



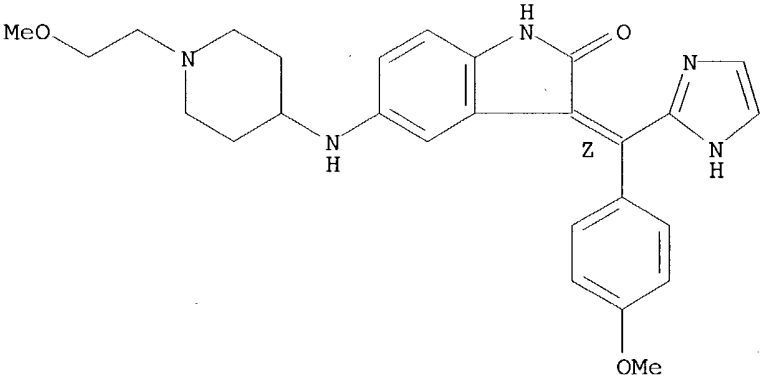
RN 705946-65-2 CAPLUS
CN 2H-Indol-2-one, 1,3-dihydro-3-(1H-imidazol-2-ylphenylmethylene)-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



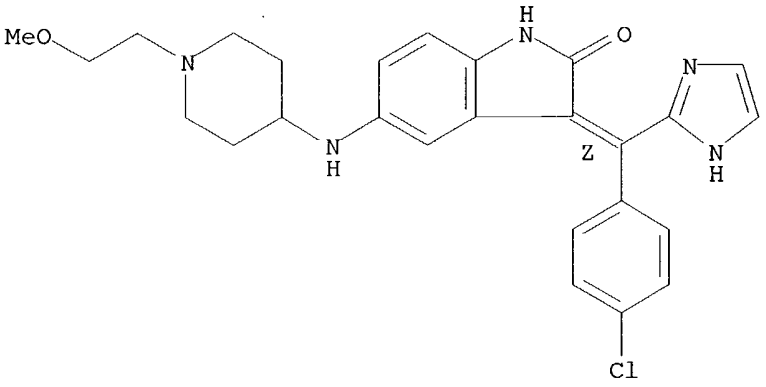
RN 705946-66-3 CAPLUS
CN 2H-Indol-2-one, 1,3-dihydro-3-[1H-imidazol-2-yl(4-methoxyphenyl)methylene]-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



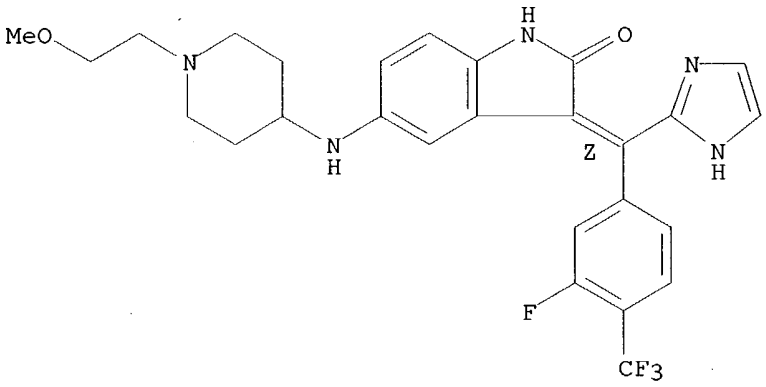
RN 705946-67-4 CAPLUS
CN 2H-Indol-2-one, 3-[(4-chlorophenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



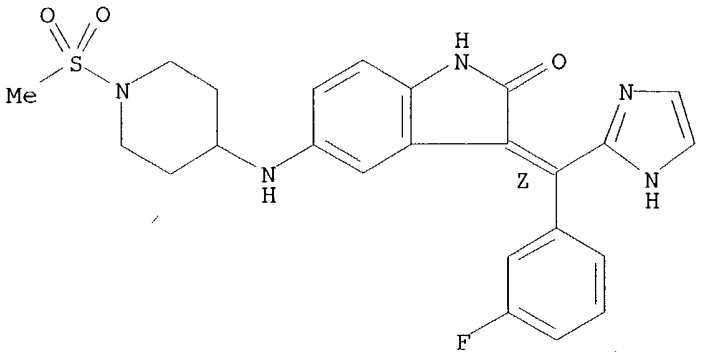
RN 705946-68-5 CAPLUS
CN 2H-Indol-2-one, 3-[[3-fluoro-4-(trifluoromethyl)phenyl]-1H-imidazol-2-ylmethylene]-1,3-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



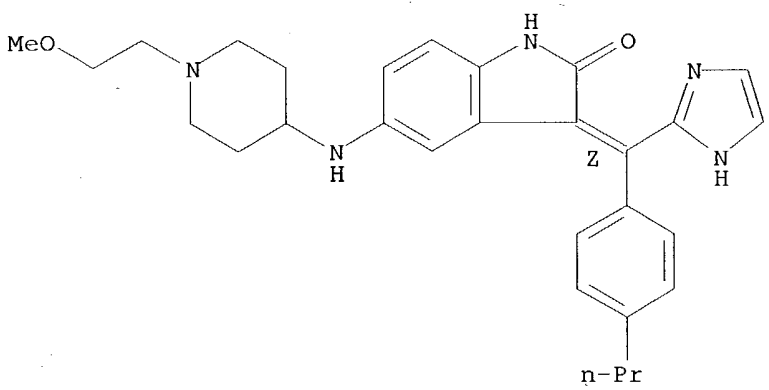
RN 705946-69-6 CAPLUS
CN 4-Piperidinamine, N-[(3Z)-3-[(3-fluorophenyl)-1H-imidazol-2-ylmethylene]-2,3-dihydro-2-oxo-1H-indol-5-yl]-1-(methylsulfonyl)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



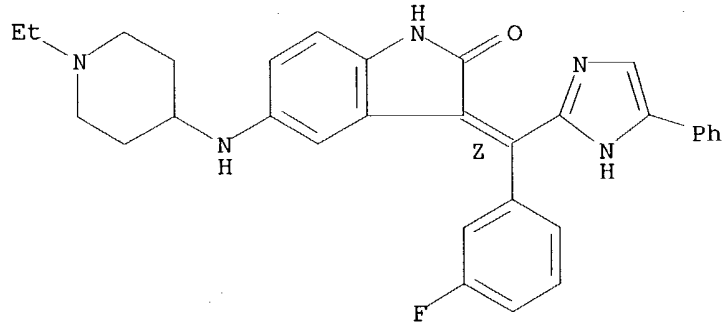
RN 705946-70-9 CAPLUS
CN 2H-Indol-2-one, 1,3-dihydro-3-[1H-imidazol-2-yl(4-propylphenyl)methylene]-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



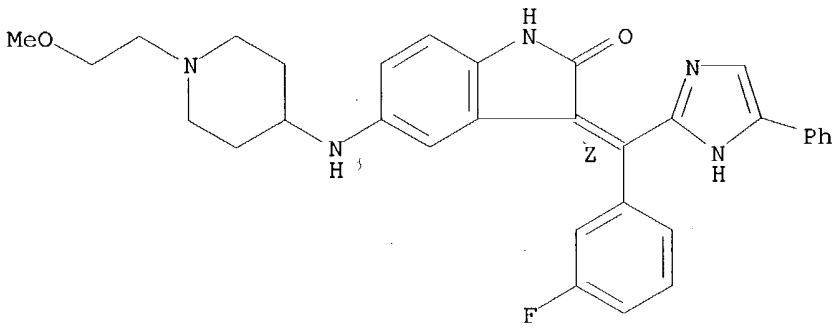
RN 705946-71-0 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[(3-fluorophenyl)(4-phenyl-1H-imidazol-2-yl)methylene]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



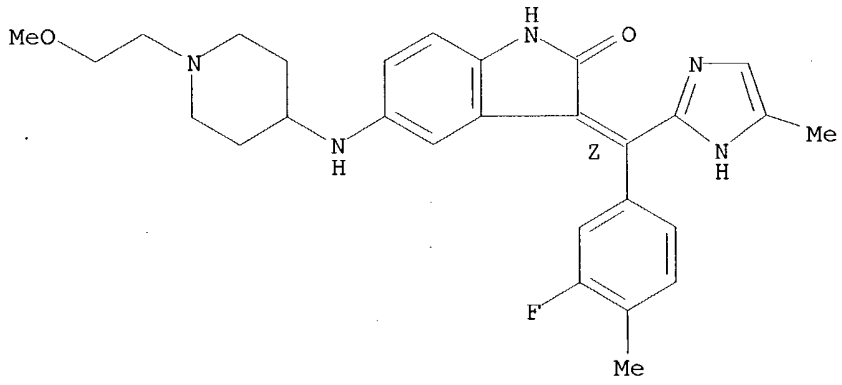
RN 705946-72-1 CAPLUS
CN 2H-Indol-2-one, 3-[(3-fluorophenyl)(4-phenyl-1H-imidazol-2-yl)methylene]-1,3-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



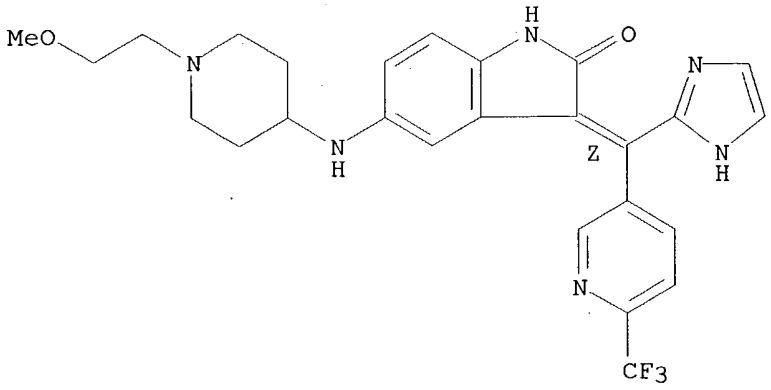
RN 705946-73-2 CAPLUS
CN 2H-Indol-2-one, 3-[(3-fluoro-4-methylphenyl)(4-methyl-1H-imidazol-2-yl)methylene]-1,3-dihydro-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



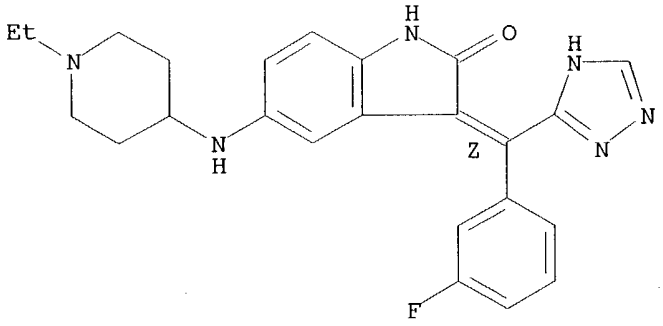
RN 705946-74-3 CAPLUS
CN 2H-Indol-2-one, 1,3-dihydro-3-[1H-imidazol-2-yl[6-(trifluoromethyl)-3-pyridinyl]methylene]-5-[[1-(2-methoxyethyl)-4-piperidinyl]amino]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



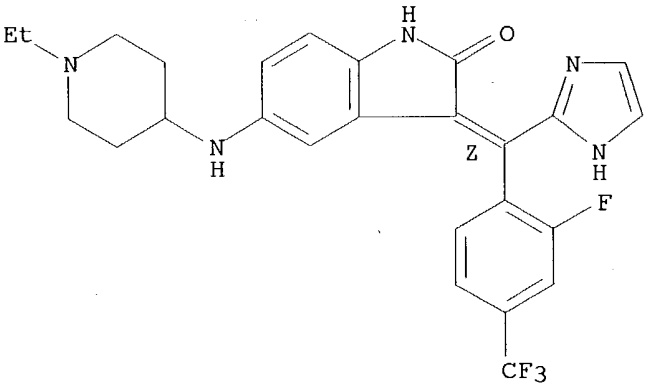
RN 705946-75-4 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[(3-fluorophenyl)-1H-1,2,4-triazol-3-ylmethylene]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



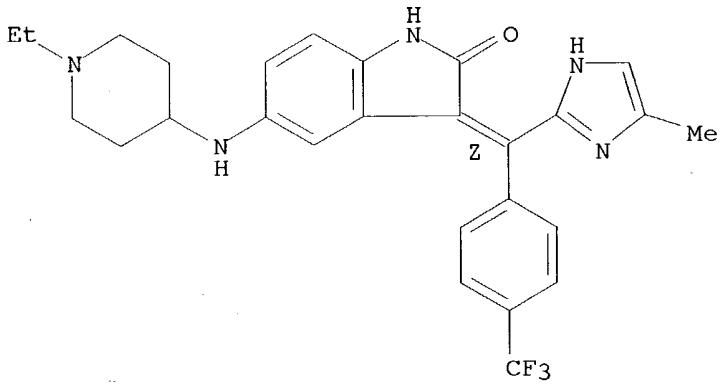
RN 705946-76-5 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[[2-fluoro-4-(trifluoromethyl)phenyl]-1H-imidazol-2-ylmethylene]-1,3-dihydro-, (3Z)-(9CI) (CA INDEX NAME)

Double bond geometry as shown.



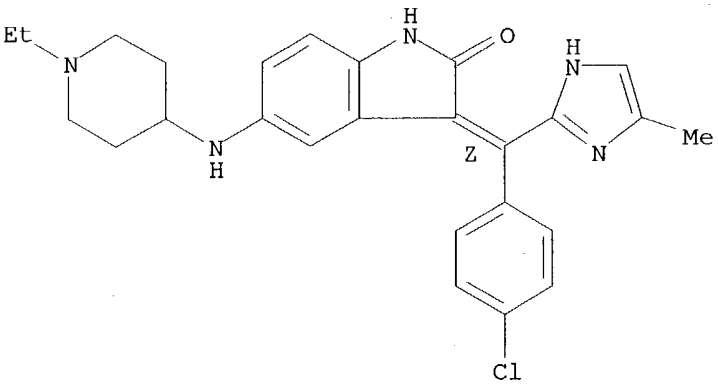
RN 705946-77-6 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-3-[(4-methyl-1H-imidazol-2-yl)[4-(trifluoromethyl)phenyl]methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



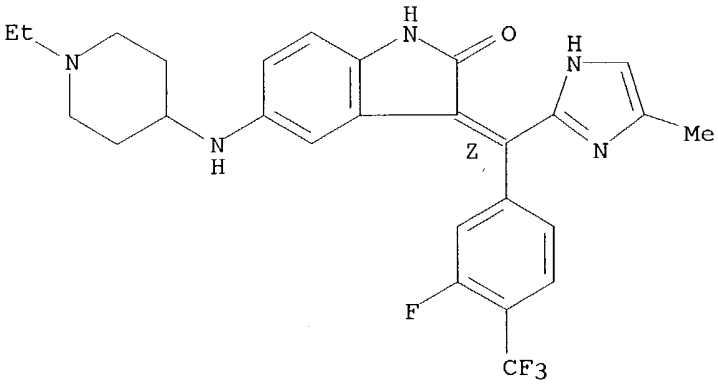
RN 705946-78-7 CAPLUS
CN 2H-Indol-2-one, 3-[(4-chlorophenyl)(4-methyl-1H-imidazol-2-yl)methylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



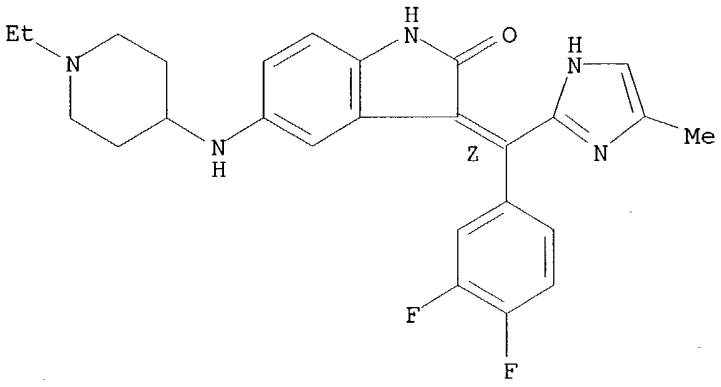
RN 705946-79-8 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[[3-fluoro-4-(trifluoromethyl)phenyl](4-methyl-1H-imidazol-2-yl)methylene]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



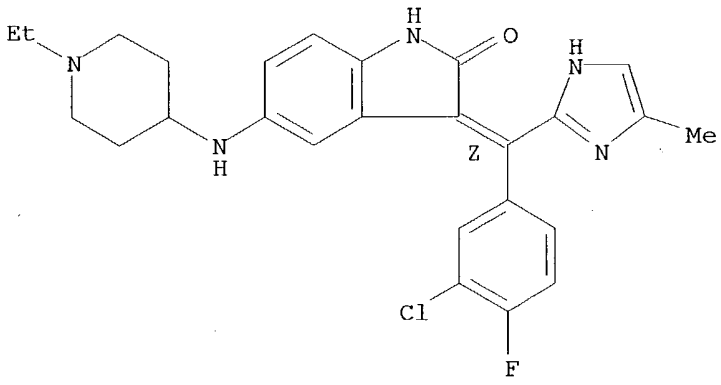
RN 705946-80-1 CAPLUS
CN 2H-Indol-2-one, 3-[(3,4-difluorophenyl)(4-methyl-1H-imidazol-2-yl)methylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



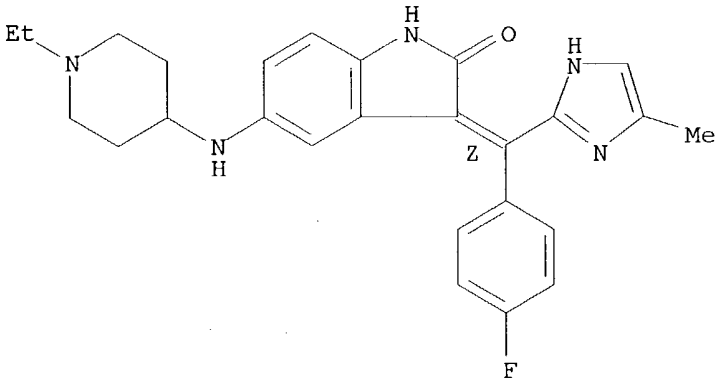
RN 705946-81-2 CAPLUS
CN 2H-Indol-2-one, 3-[(3-chloro-4-fluorophenyl)(4-methyl-1H-imidazol-2-yl)methylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



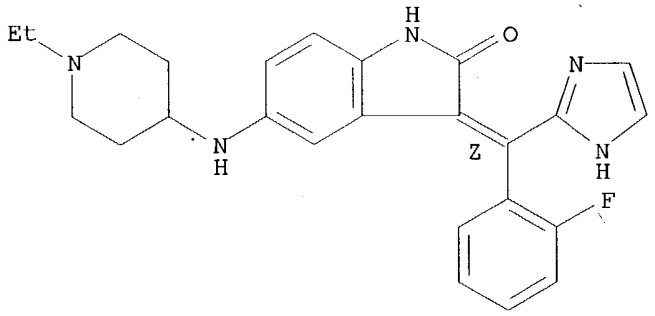
RN 705946-82-3 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[(4-fluorophenyl)(4-methyl-1H-imidazol-2-yl)methylene]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



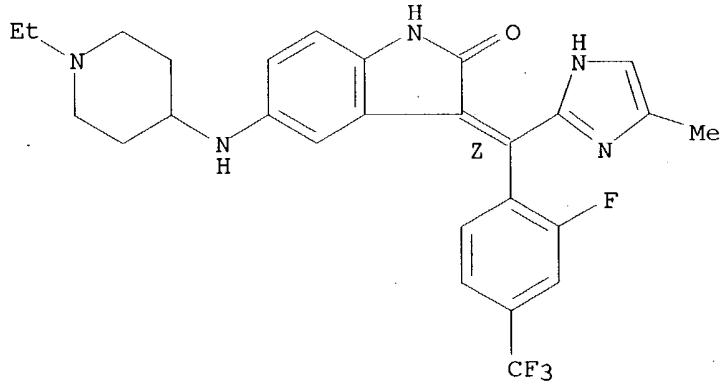
RN 705946-83-4 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[(2-fluorophenyl)-1H-imidazol-2-ylmethylene]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



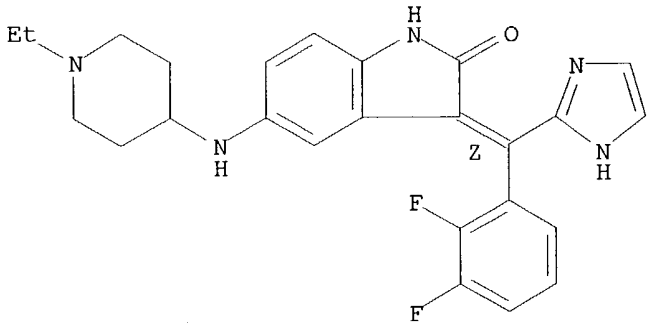
RN 705946-84-5 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[[2-fluoro-4-(trifluoromethyl)phenyl](4-methyl-1H-imidazol-2-yl)methylene]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



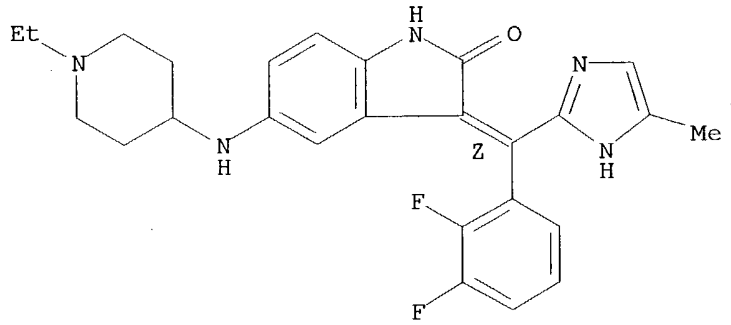
RN 705946-85-6 CAPLUS
CN 2H-Indol-2-one, 3-[(2,3-difluorophenyl)-1H-imidazol-2-ylmethylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



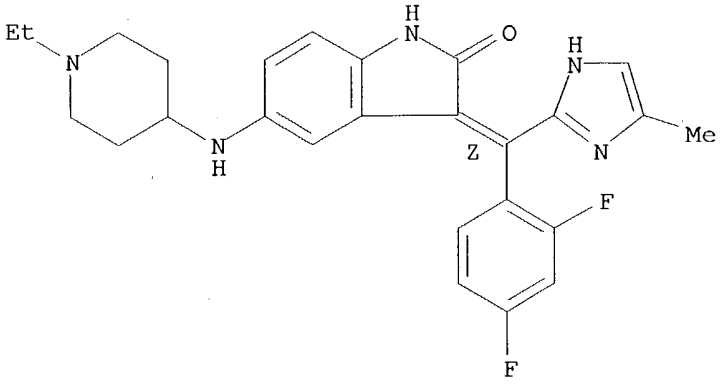
RN 705946-86-7 CAPLUS
CN 2H-Indol-2-one, 3-[(2,3-difluorophenyl)(4-methyl-1H-imidazol-2-yl)methylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



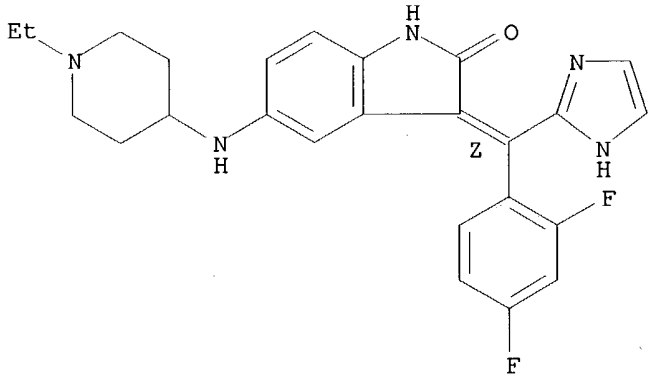
RN 705946-87-8 CAPLUS
CN 2H-Indol-2-one, 3-[(2,4-difluorophenyl)(4-methyl-1H-imidazol-2-yl)methylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



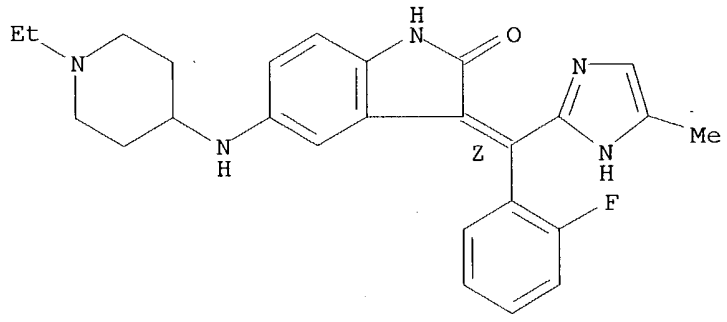
RN 705946-88-9 CAPLUS
CN 2H-Indol-2-one, 3-[(2,4-difluorophenyl)-1H-imidazol-2-ylmethylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



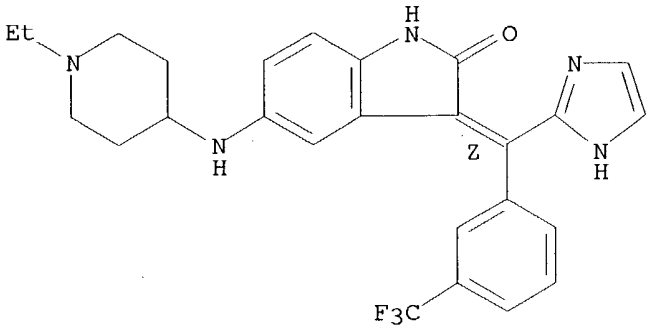
RN 705946-89-0 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[(2-fluorophenyl)(4-methyl-1H-imidazol-2-yl)methylene]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



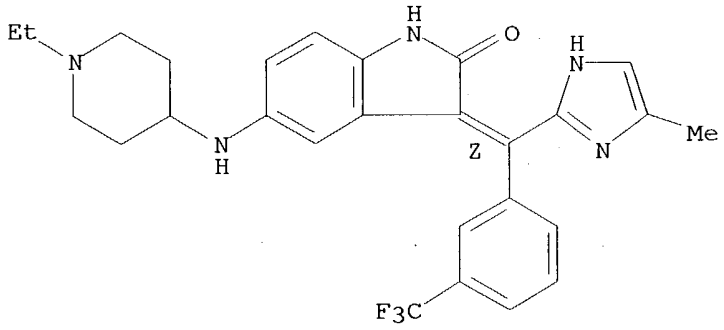
RN 705946-90-3 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-3-[1H-imidazol-2-yl[3-(trifluoromethyl)phenyl]methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



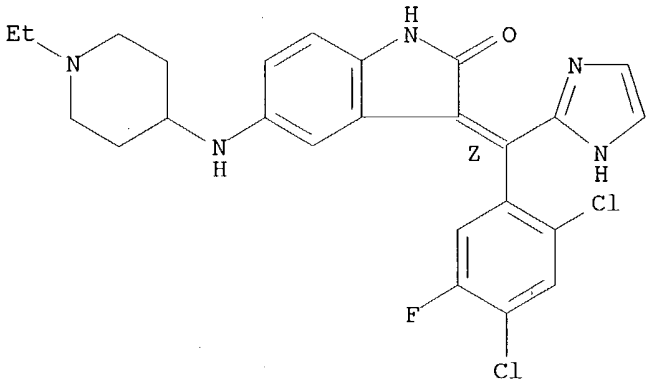
RN 705946-91-4 CAPLUS
CN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-3-[(4-methyl-1H-imidazol-2-yl)[3-(trifluoromethyl)phenyl]methylene]-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



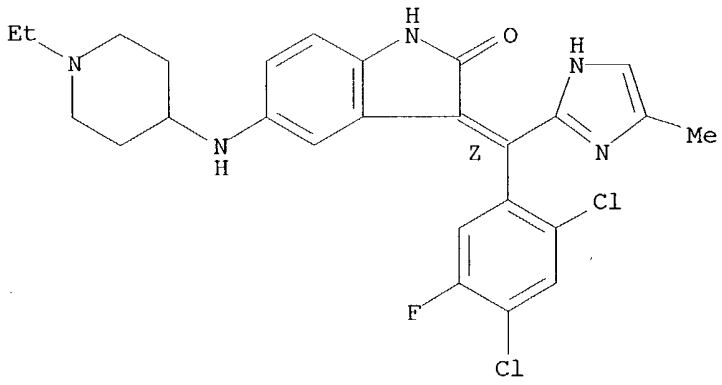
RN 705946-92-5 CAPLUS
CN 2H-Indol-2-one, 3-[(2,4-dichloro-5-fluorophenyl)-1H-imidazol-2-ylmethylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



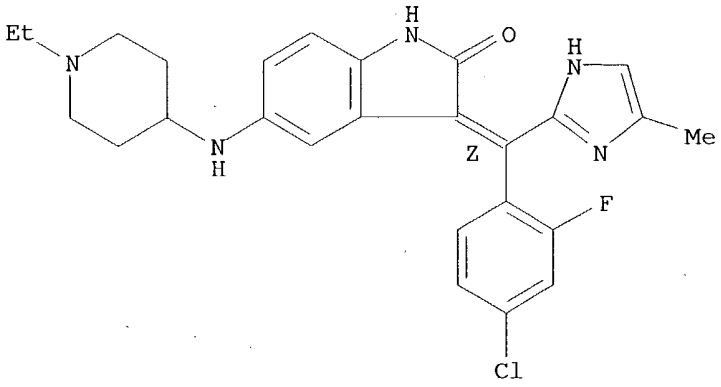
RN 705946-93-6 CAPLUS
CN 2H-Indol-2-one, 3-[(2,4-dichloro-5-fluorophenyl)(4-methyl-1H-imidazol-2-yl)methylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 705946-94-7 CAPLUS
CN 2H-Indol-2-one, 3-[(4-chloro-2-fluorophenyl)(4-methyl-1H-imidazol-2-yl)methylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.


Et
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Me
Z
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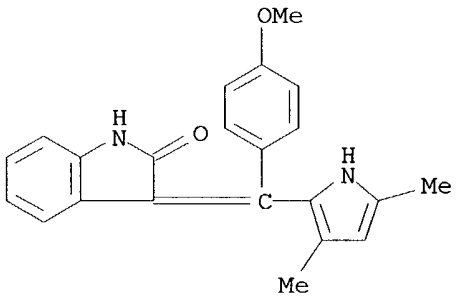
Cl

V. Balasubramanian

L5 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2004 ACS on STN
AN 2000:688215 CAPLUS
DN 133:252306
TI Preparation of indolinones as protein kinase inhibitors.
IN Tang, Peng Cho; Sun, Li; Mcmahon, Gerald; Miller, Todd Anthony; Shirazian, Shahrzad; Wei, Chung Chen; Harris, G. Davis; Xiaoyuan, Li; Liang, Congxin
PA Sugen, Inc., USA
SO PCT Int. Appl., 245 pp.
CODEN: PIXXD2
DT Patent
LA English
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2000056709	A1	20000928	WO 2000-US7704	20000322
	W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW: GH, GM, KE, LS, MW, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
	EP 1165513	A1	20020102	EP 2000-916622	20000322
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO				
	JP 2002540096	T2	20021126	JP 2000-606571	20000322
	US 6689806	B1	20040210	US 2000-534405	20000322
PRAI	US 1999-125945P	P	19990324		
	US 1999-127863P	P	19990405		
	US 1999-131192P	P	19990426		
	US 1999-132243P	P	19990503		
	WO 2000-US7704	W	20000322		

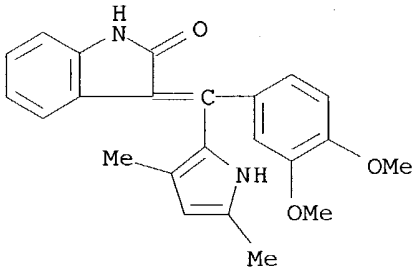
OS MARPAT 133:252306
IT **295799-89-2P 295799-90-5P**
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of indolinones as protein kinase inhibitors)
RN 295799-89-2 CAPLUS
CN 2H-Indol-2-one, 3-[(3,5-dimethyl-1H-pyrrol-2-yl)(4-methoxyphenyl)methylene]-1,3-dihydro- (9CI) (CA INDEX NAME)



RN 295799-90-5 CAPLUS

V. Balasubramanian

CN 2H-Indol-2-one, 3-[(3,4-dimethoxyphenyl)(3,5-dimethyl-1H-pyrrol-2-yl)methylene]-1,3-dihydro- (9CI) (CA INDEX NAME)



RE.CNT 31 THERE ARE 31 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

V. Balasubramanian

=> log y

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	6.98	162.61

STN INTERNATIONAL LOGOFF AT 13:24:06 ON 24 AUG 2004

V. Balasubramanian

Welcome to STN International! Enter x:x

LOGINID:ssspta1611bxv

PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * * * * * * Welcome to STN International * * * * * * * * * *

NEWS 1		Web Page URLs for STN Seminar Schedule - N. America
NEWS 2		"Ask CAS" for self-help around the clock
NEWS 3	May 12	EXTEND option available in structure searching
NEWS 4	May 12	Polymer links for the POLYLINK command completed in REGISTRY
NEWS 5	May 27	New UPM (Update Code Maximum) field for more efficient patent SDIs in CAplus
NEWS 6	May 27	CAplus super roles and document types searchable in REGISTRY
NEWS 7	Jun 28	Additional enzyme-catalyzed reactions added to CASREACT
NEWS 8	Jun 28	ANTE, AQUALINE, BIOENG, CIVILENG, ENVIROENG, MECHENG, and WATER from CSA now available on STN(R)
NEWS 9	Jul 12	BEILSTEIN enhanced with new display and select options, resulting in a closer connection to BABS
NEWS 10	Jul 30	BEILSTEIN on STN workshop to be held August 24 in conjunction with the 228th ACS National Meeting
NEWS 11	AUG 02	IFIPAT/IFIUDB/IFICDB reloaded with new search and display fields
NEWS 12	AUG 02	CAplus and CA patent records enhanced with European and Japan Patent Office Classifications
NEWS 13	AUG 02	STN User Update to be held August 22 in conjunction with the 228th ACS National Meeting
NEWS 14	AUG 02	The Analysis Edition of STN Express with Discover! (Version 7.01 for Windows) now available
NEWS 15	AUG 04	Pricing for the Save Answers for SciFinder Wizard within STN Express with Discover! will change September 1, 2004

NEWS EXPRESS	JULY 30 CURRENT WINDOWS VERSION IS V7.01, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 11 AUGUST 2004
NEWS HOURS	STN Operating Hours Plus Help Desk Availability
NEWS INTER	General Internet Information
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NEWS PHONE	Direct Dial and Telecommunication Network Access to STN
NEWS WWW	CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

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* * * * * * * * * * * * * STN Columbus * * * * * * * * * * * * * * * *

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FILE 'HOME' ENTERED AT 14:09:21 ON 24 AUG 2004

=> file reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'REGISTRY' ENTERED AT 14:09:31 ON 24 AUG 2004
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 23 AUG 2004 HIGHEST RN 731771-88-3
DICTIONARY FILE UPDATES: 23 AUG 2004 HIGHEST RN 731771-88-3

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

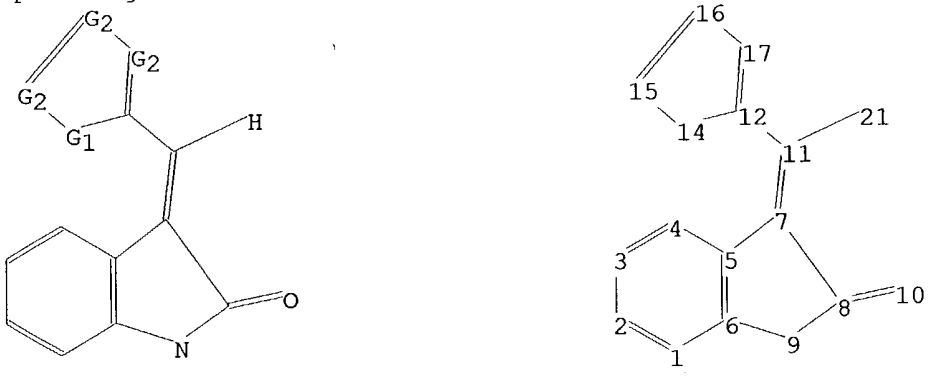
Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:
http://www.cas.org/ONLINE/DBSS/registryss.html

=> ....Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=>
Uploading C:\STNEXP4\QUERIES\10725079-7.str



chain nodes :
10 11 21
ring nodes :
1 2 3 4 5 6 7 8 9 12 14 15 16 17
chain bonds :
7-11 8-10 11-12 11-21

V. Balasubramanian

ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 12-14 12-17 14-15 15-16 16-17

exact/norm bonds :
5-7 6-9 7-8 7-11 8-9 8-10 11-12 11-21 12-14 12-17 14-15 15-16 16-17
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6
isolated ring systems :
containing 1 : 12 :

G1:O,S,N

G2:C,O,S,N

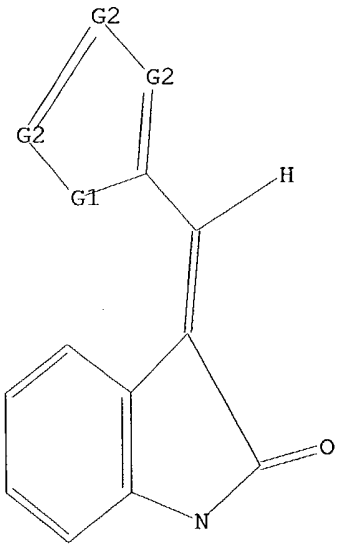
Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS
11:CLASS 12:Atom 14:Atom 15:Atom 16:Atom 17:Atom 21:CLASS

L1 STRUCTURE UPLOADED

=> que L1

L2 QUE L1

=> d l1
L1 HAS NO ANSWERS
L1 STR



G1 O,S,N
G2 C,O,S,N

Structure attributes must be viewed using STN Express query preparation.

```
=> s l1 sss sam
SAMPLE SEARCH INITIATED 14:09:53 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -     634 TO ITERATE

100.0% PROCESSED      634 ITERATIONS                                 50 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:        11170 TO    14190
PROJECTED ANSWERS:            1934 TO     3306

L3            50 SEA SSS SAM L1

=> d scan
```

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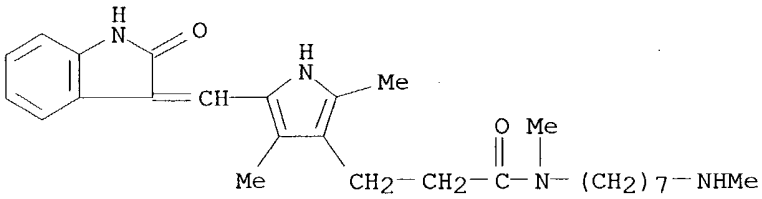
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-propanamide, 5-[(1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-N,2,4-trimethyl-N-[7-(methylamino)heptyl]- (9CI)
MF C27 H38 N4 O2
CI COM



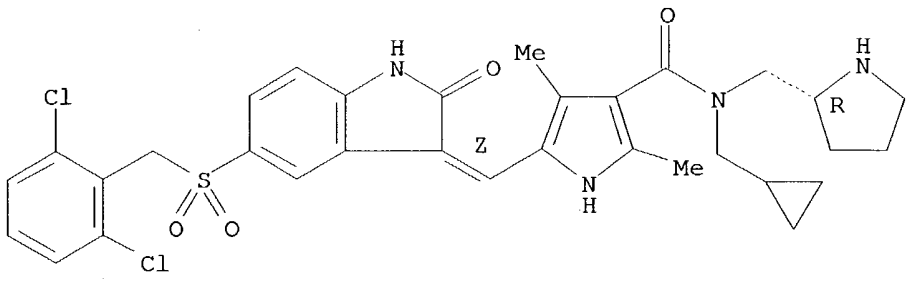
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):49

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-carboxamide, N-(cyclopropylmethyl)-5-[(Z)-[5-[[(2,6-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-N-[(2R)-2-pyrrolidinylmethyl]- (9CI)
MF C32 H34 Cl2 N4 O4 S

Absolute stereochemistry.
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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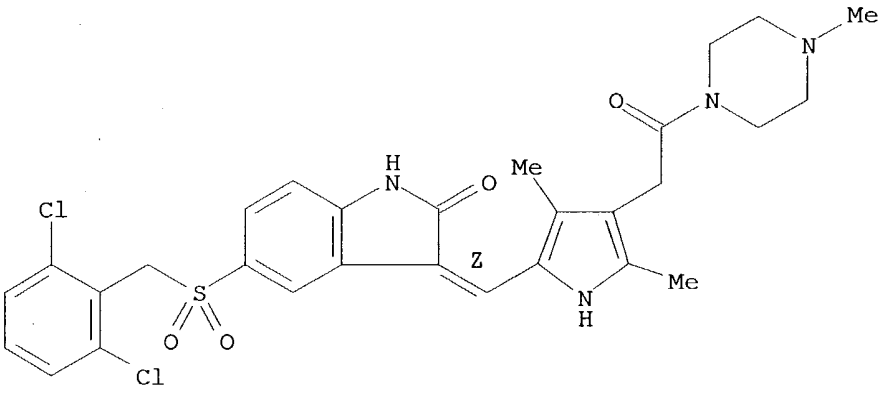
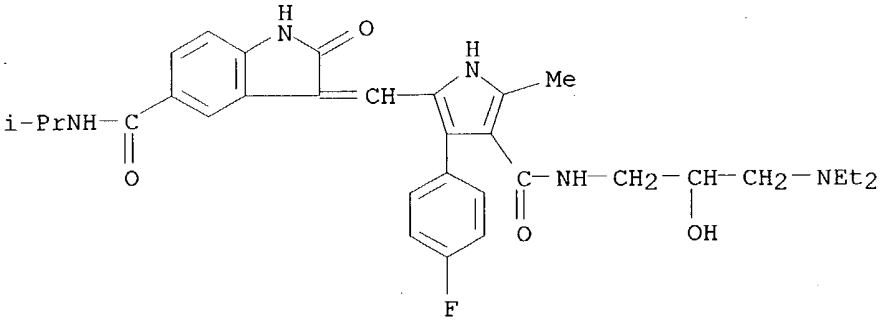
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperazine, 1-[[5-[(Z)-[5-[[(2,6-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]acetyl]-4-methyl- (9CI)
MF C29 H30 Cl2 N4 O4 S

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

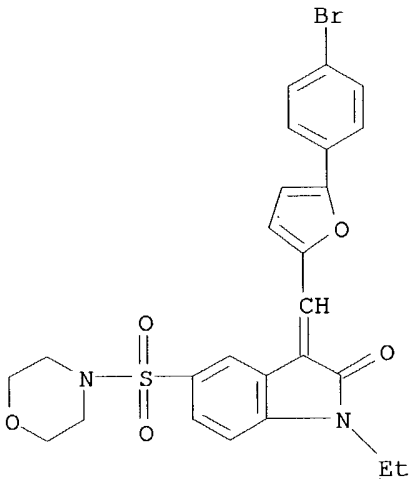
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Indole-5-carboxamide, 3-[[4-[[[3-(diethylamino)-2-hydroxypropyl]amino]carbonyl]-3-(4-fluorophenyl)-5-methyl-1H-pyrrol-2-yl]methylene]-2,3-dihydro-N-(1-methylethyl)-2-oxo- (9CI)
MF C32 H38 F N5 O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Morpholine, 4-[[3-[[5-(4-bromophenyl)-2-furanyl]methylene]-1-ethyl-2,3-dihydro-2-oxo-1H-indol-5-yl]sulfonyl]- (9CI)
MF C25 H23 Br N2 O5 S

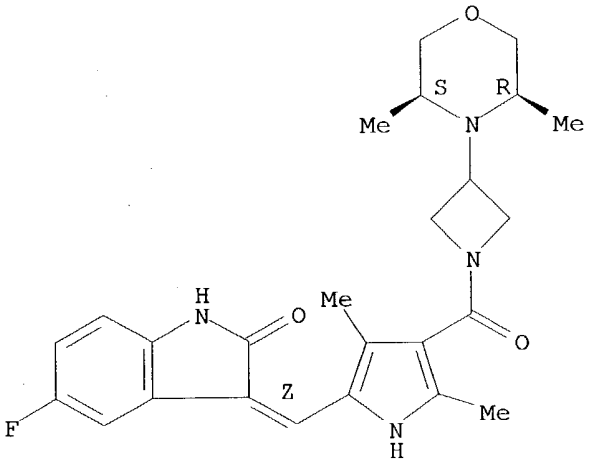


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Azetidine, 3-[(3R,5S)-3,5-dimethyl-4-morpholinyl]-1-[[5-[(Z)-(5-fluoro-1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-, rel- (9CI)
MF C25 H29 F N4 O3

Relative stereochemistry.
Double bond geometry as shown.

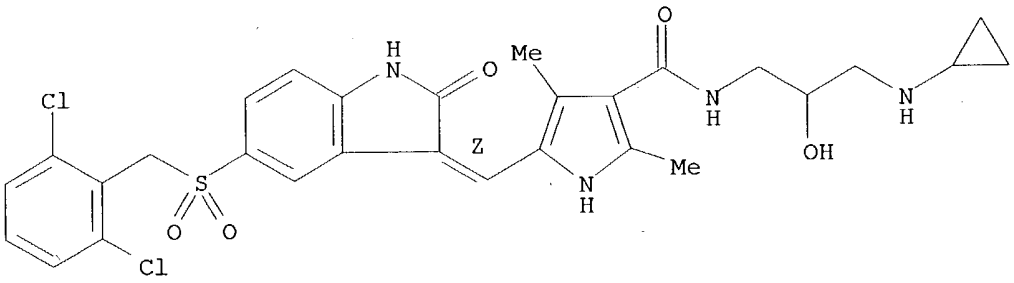


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-carboxamide, N-[3-(cyclopropylamino)-2-hydroxypropyl]-5-[(Z)-[5-[[(2,6-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl- (9CI)
MF C29 H30 Cl2 N4 O5 S

Double bond geometry as shown.

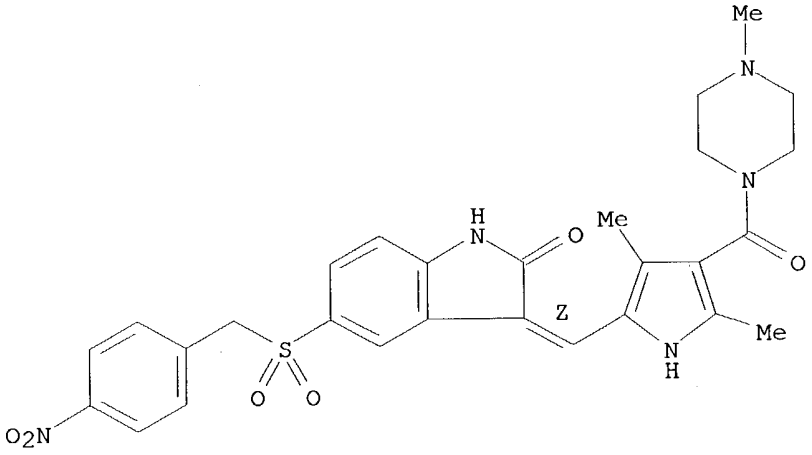


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperazine, 1-[[5-[(Z)-[1,2-dihydro-5-[[(4-nitrophenyl)methyl]sulfonyl]-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-4-methyl- (9CI)
MF C28 H29 N5 O6 S

Double bond geometry as shown.

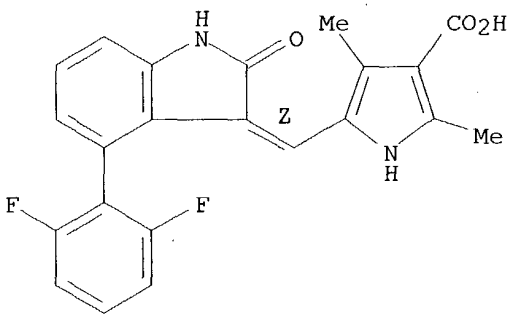


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-carboxylic acid, 5-[(Z)-[4-(2,6-difluorophenyl)-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl- (9CI)
MF C22 H16 F2 N2 O3

Double bond geometry as shown.

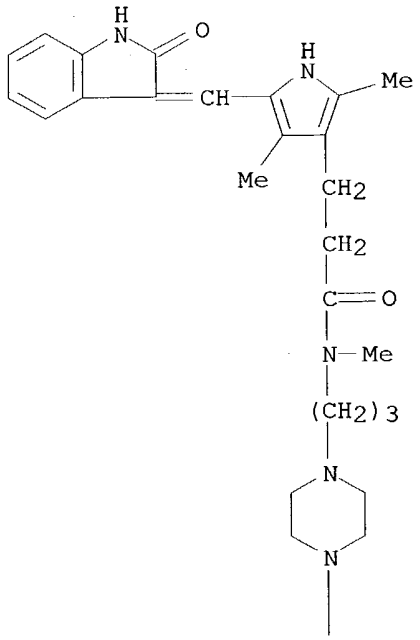


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

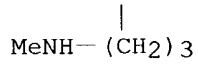
V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-propanamide, 5-[(1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-N,2,4-trimethyl-N-[3-[4-[3-(methylamino)propyl]-1-piperazinyl]propyl]-(9CI)
MF C30 H44 N6 O2
CI COM

PAGE 1-A



PAGE 2-A

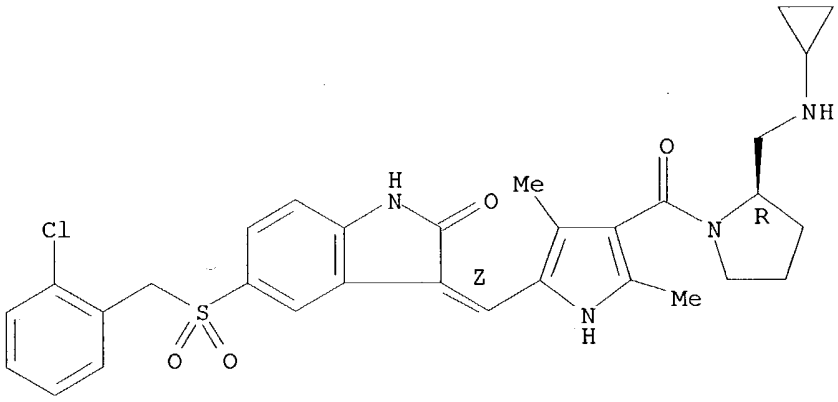


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2-Pyrrolidinemethanamine, 1-[[5-[(Z)-[5-[[(2-chlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-N-cyclopropyl-, (2R)- (9CI)
MF C31 H33 Cl N4 O4 S

Absolute stereochemistry.
Double bond geometry as shown.

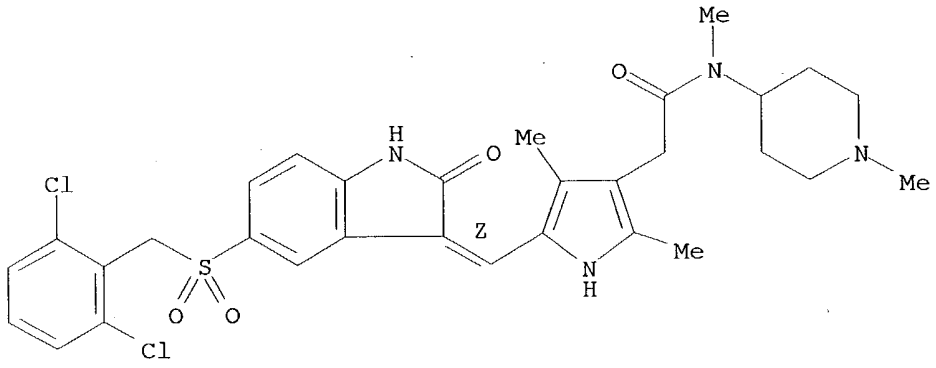


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-acetamide, 5-[(Z)-[5-[[(2,6-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-N,2,4-trimethyl-N-(1-methyl-4-piperidinyl)- (9CI)
MF C31 H34 Cl2 N4 O4 S

Double bond geometry as shown.

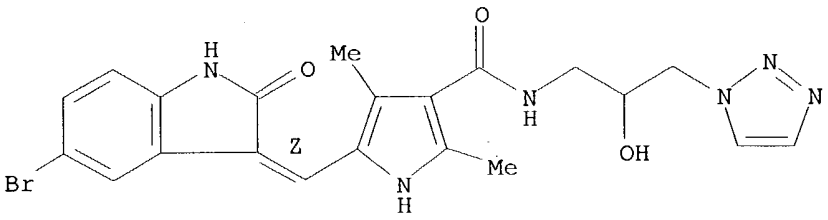


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-carboxamide, 5-[(Z)-(5-bromo-1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-N-[2-hydroxy-3-(1H-1,2,3-triazol-1-yl)propyl]-2,4-dimethyl- (9CI)
MF C21 H21 Br N6 O3

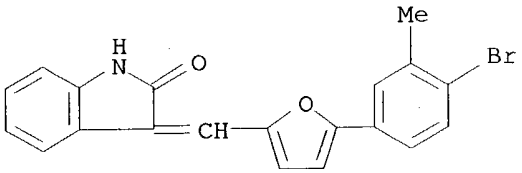
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 3-[[5-(4-bromo-3-methylphenyl)-2-furanyl]methylene]-1,3-dihydro- (9CI)
MF C20 H14 Br N O2

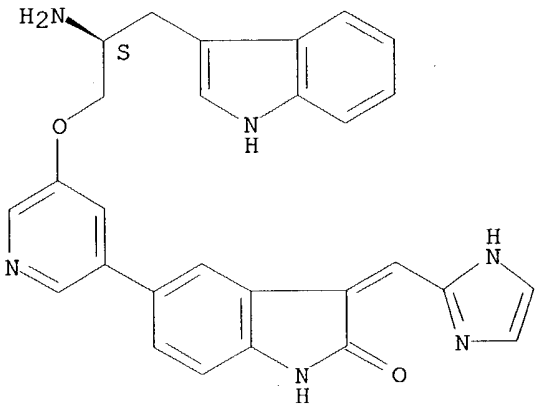


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 5-[5-[(2S)-2-amino-3-(1H-indol-3-yl)propoxy]-3-pyridinyl]-1,3-dihydro-3-(1H-imidazol-2-ylmethylene)- (9CI)
MF C28 H24 N6 O2
CI COM

Absolute stereochemistry.
Double bond geometry unknown.

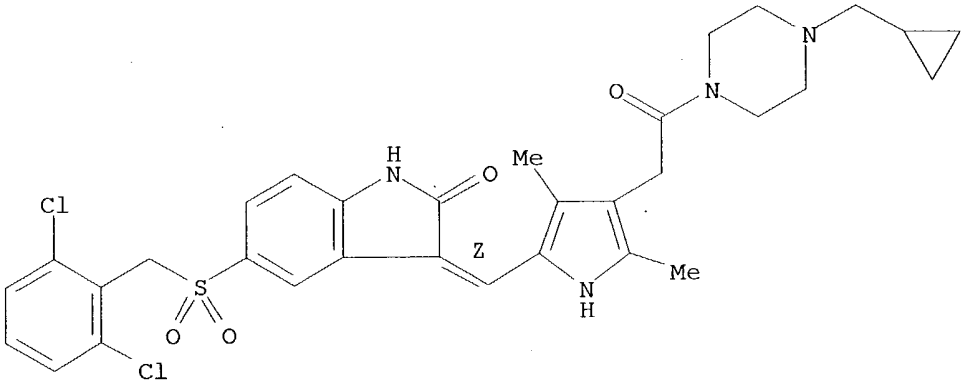


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperazine, 1-(cyclopropylmethyl)-4-[[5-[(Z)-[5-[[(2,6-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]acetyl]- (9CI)
MF C32 H34 Cl2 N4 O4 S

Double bond geometry as shown.

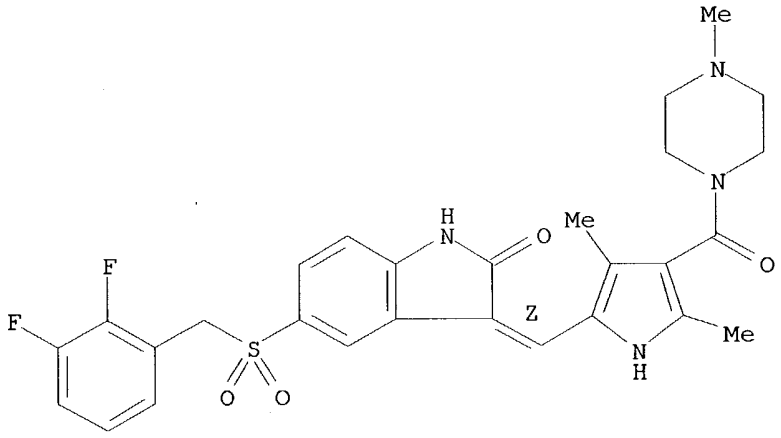


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperazine, 1-[[5-[(Z)-[5-[[(2,3-difluorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-4-methyl- (9CI)
MF C28 H28 F2 N4 O4 S

Double bond geometry as shown.

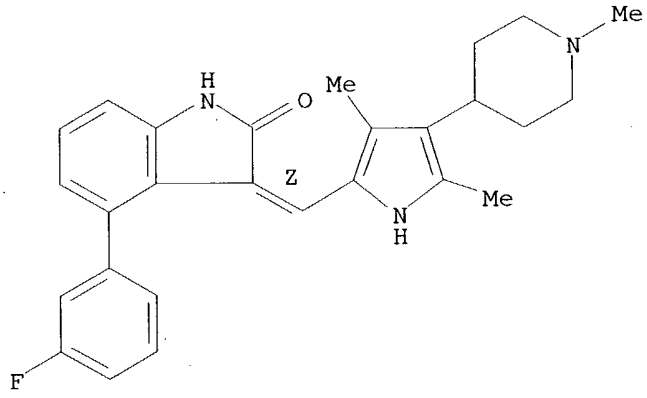


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 3-[[3,5-dimethyl-4-(1-methyl-4-piperidinyl)-1H-pyrrol-2-yl]methylene]-4-(3-fluorophenyl)-1,3-dihydro-, (3Z)- (9CI)
MF C27 H28 F N3 O

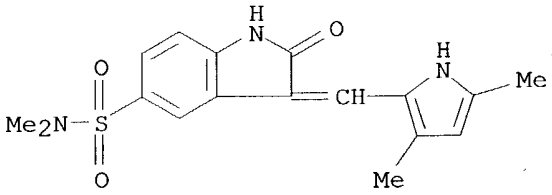
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Indole-5-sulfonamide, 3-[(3,5-dimethyl-1H-pyrrol-2-yl)methylene]-2,3-dihydro-N,N-dimethyl-2-oxo- (9CI)
MF C17 H19 N3 O3 S

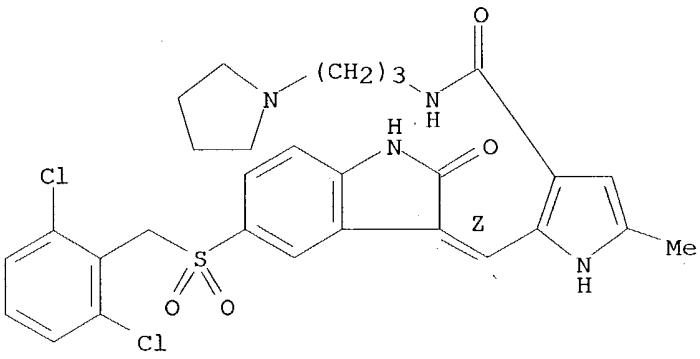


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-carboxamide, 2-[(Z)-[5-[[(2,6-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-5-methyl-N-[3-(1-pyrrolidinyl)propyl]- (9CI)
MF C29 H30 Cl2 N4 O4 S

Double bond geometry as shown.

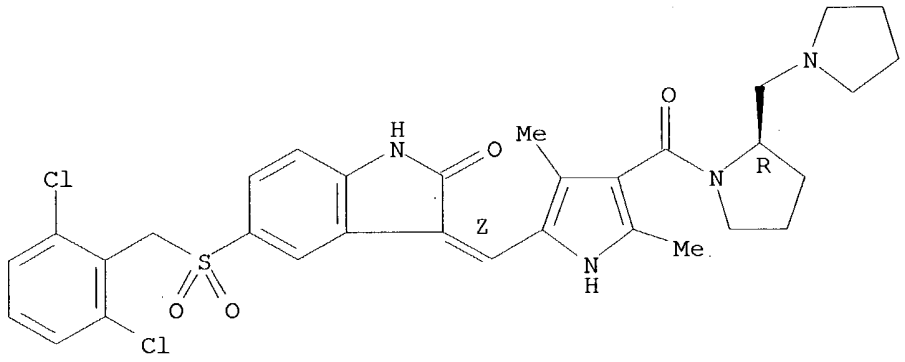


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Pyrrolidine, 1-[[5-[(Z)-[5-[[(2,6-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-2-(1-pyrrolidinylmethyl)-, (2R)- (9CI)
MF C32 H34 Cl2 N4 O4 S

Absolute stereochemistry.
Double bond geometry as shown.

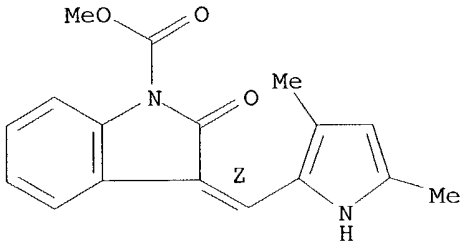


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Indole-1-carboxylic acid, 3-[(3,5-dimethyl-1H-pyrrol-2-yl)methylene]-2,3-dihydro-2-oxo-, methyl ester, (3Z)- (9CI)
MF C17 H16 N2 O3

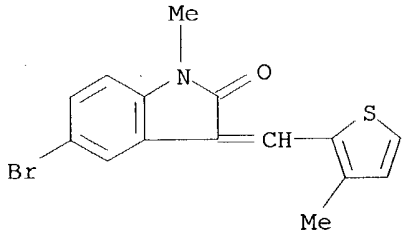
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 5-bromo-1,3-dihydro-1-methyl-3-[(3-methyl-2-thienyl)methylene]- (9CI)
MF C15 H12 Br N O S

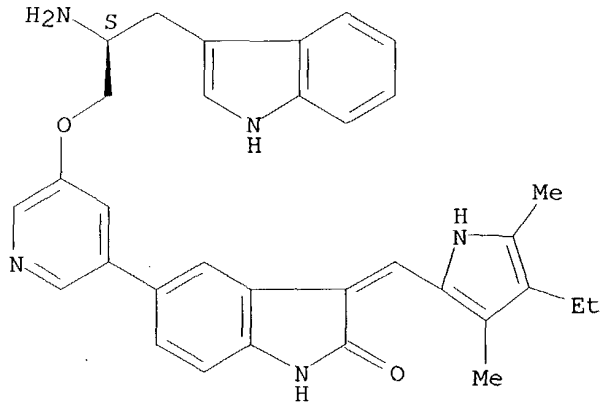


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 5-[5-[(2S)-2-amino-3-(1H-indol-3-yl)propoxy]-3-pyridinyl]-3-[(4-ethyl-3,5-dimethyl-1H-pyrrol-2-yl)methylene]-1,3-dihydro- (9CI)
MF C33 H33 N5 O2
CI COM

Absolute stereochemistry.
Double bond geometry unknown.

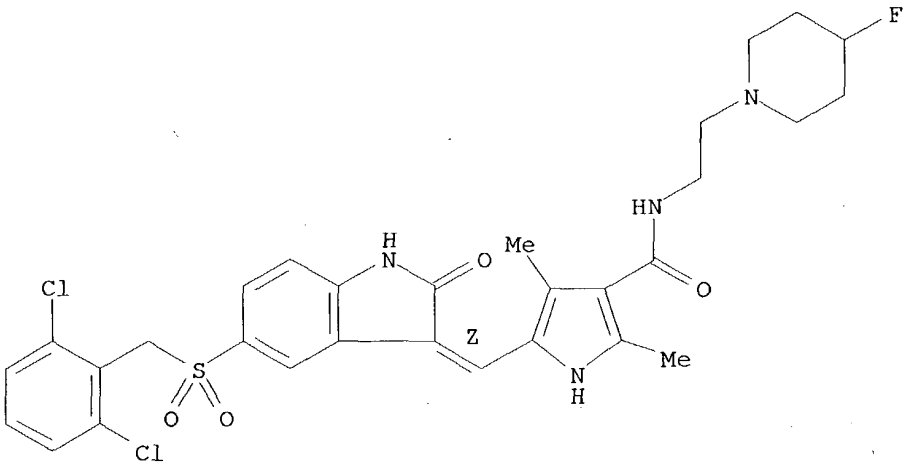


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-carboxamide, 5-[(Z)-[5-[[(2,6-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-N-[2-(4-fluoro-1-piperidinyl)ethyl]-2,4-dimethyl- (9CI)
MF C30 H31 Cl2 F N4 O4 S

Double bond geometry as shown.

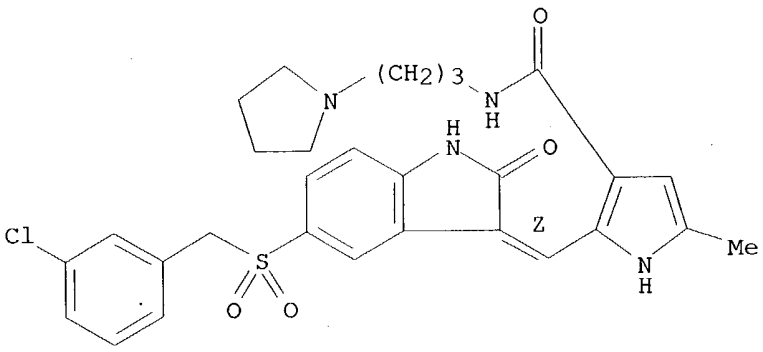


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-carboxamide, 2-[(Z)-[5-[[(3-chlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-5-methyl-N-[3-(1-pyrrolidinyl)propyl]- (9CI)
MF C29 H31 Cl N4 O4 S

Double bond geometry as shown.

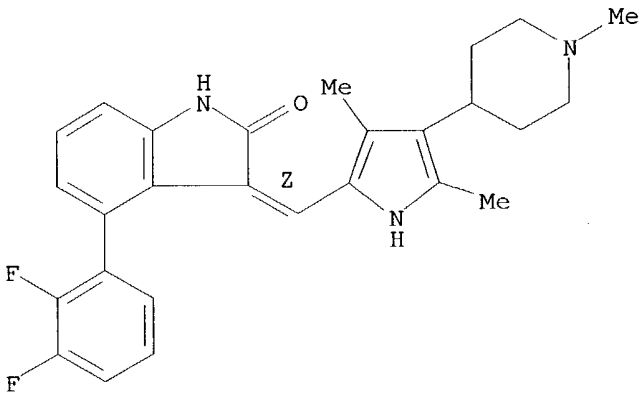


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 4-(2,3-difluorophenyl)-3-[[3,5-dimethyl-4-(1-methyl-4-piperidinyl)-1H-pyrrol-2-yl]methylene]-1,3-dihydro-, (3Z)- (9CI)
MF C27 H27 F2 N3 O

Double bond geometry as shown.

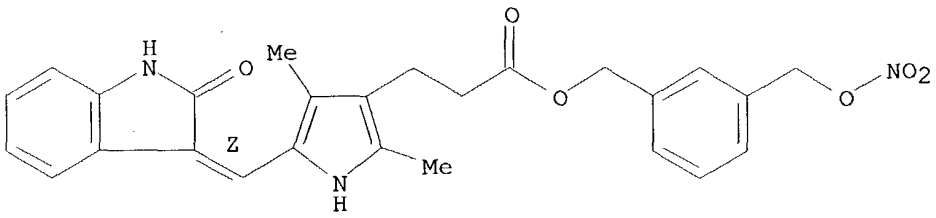


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-propanoic acid, 5-[(Z)-(1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-2,4-dimethyl-, [3-[(nitrooxy)methyl]phenyl]methyl ester (9CI)
MF C26 H25 N3 O6

Double bond geometry as shown.

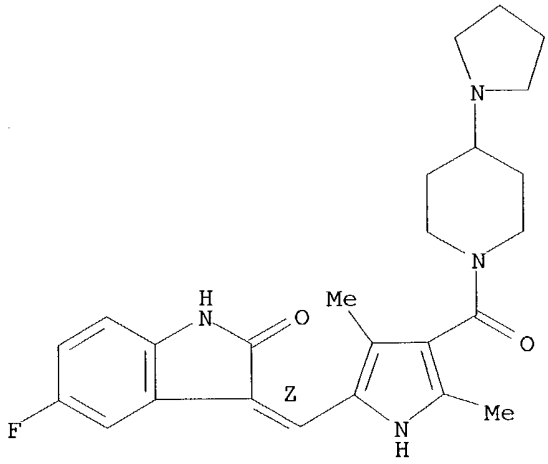


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperidine, 1-[[5-[(Z)-(5-fluoro-1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-4-(1-pyrrolidinyl)-(9CI)
MF C25 H29 F N4 O2

Double bond geometry as shown.

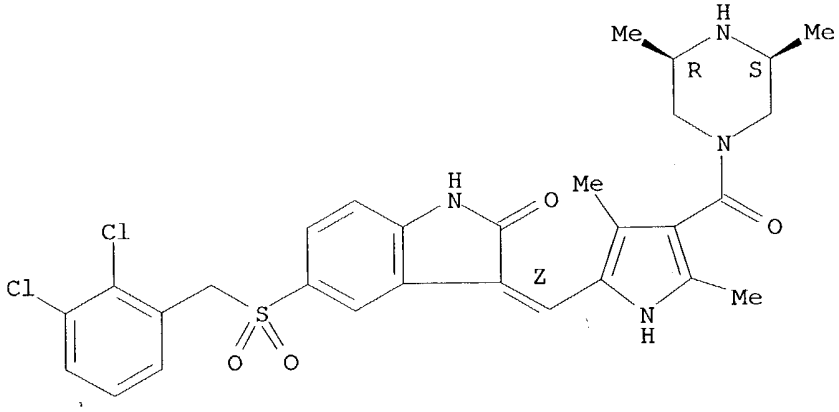


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperazine, 1-[[5-[(Z)-[5-[[(2,3-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-3,5-dimethyl-, (3R,5S)- (9CI)
MF C29 H30 Cl2 N4 O4 S

Absolute stereochemistry.
Double bond geometry as shown.

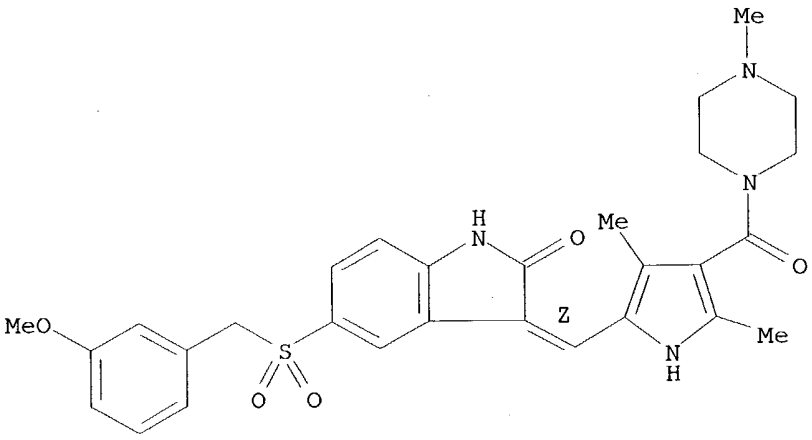


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperazine, 1-[[5-[(Z)-[1,2-dihydro-5-[[(3-methoxyphenyl)methyl]sulfonyl]-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-4-methyl- (9CI)
MF C29 H32 N4 O5 S

Double bond geometry as shown.

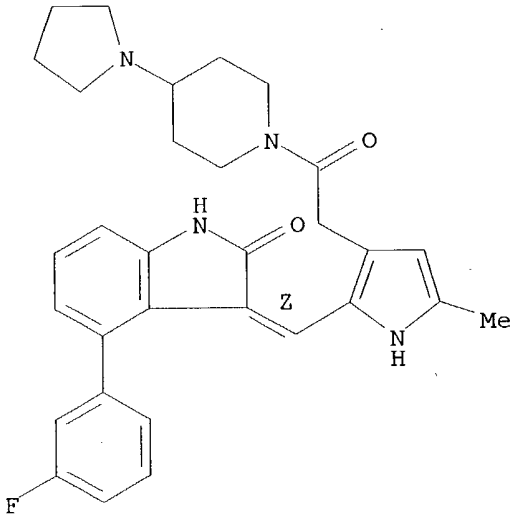


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

V. Balasubramanian

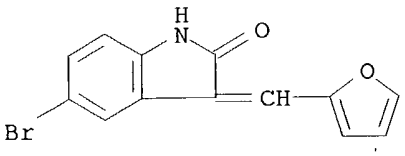
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperidine, 1-[[2-[(Z)-[4-(3-fluorophenyl)-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-5-methyl-1H-pyrrol-3-yl]acetyl]-4-(1-pyrrolidinyl)- (9CI)
MF C31 H33 F N4 O2

Double bond geometry as shown.



V. Balasubramanian

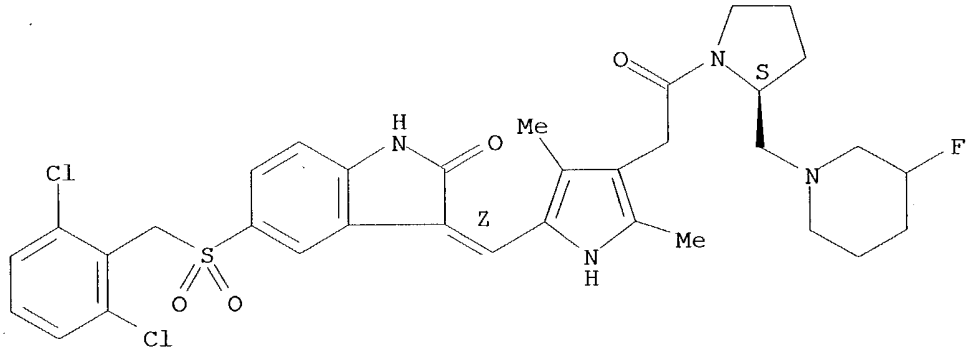
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 5-bromo-3-(2-furanylmethylene)-1,3-dihydro- (9CI)
MF C13 H8 Br N O2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Pyrrolidine, 1-[[5-[(Z)-[5-[[(2,6-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]acetyl]-2-[(3-fluoro-1-piperidinyl)methyl]-, (2S)- (9CI)
MF C34 H37 Cl2 F N4 O4 S

Absolute stereochemistry.
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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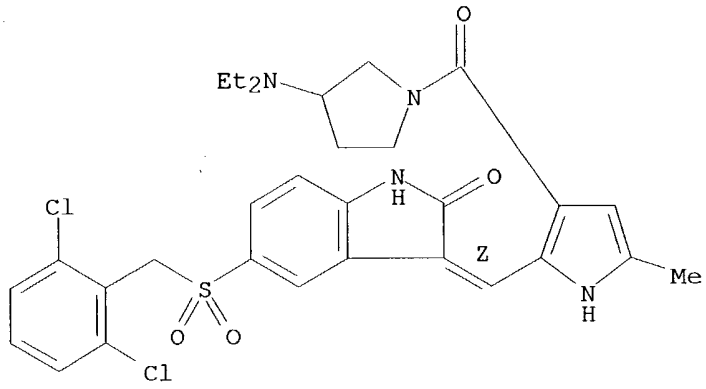
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 3-Pyrrolidinamine, 1-[[2-[(Z)-[5-[[(2,6-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-5-methyl-1H-pyrrol-3-yl]carbonyl]-N,N-diethyl- (9CI)
MF C30 H32 Cl2 N4 O4 S

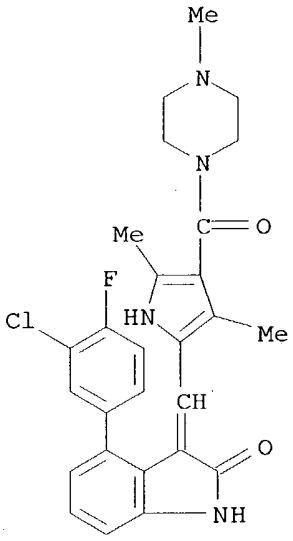
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

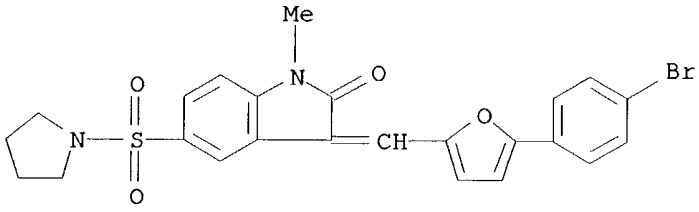
V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperazine, 1-[[5-[[4-(3-chloro-4-fluorophenyl)-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-4-methyl- (9CI)
MF C27 H26 Cl F N4 O2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Pyrrolidine, 1-[[3-[[5-(4-bromophenyl)-2-furanyl]methylene]-2,3-dihydro-1-methyl-2-oxo-1H-indol-5-yl]sulfonyl]- (9CI)
MF C24 H21 Br N2 O4 S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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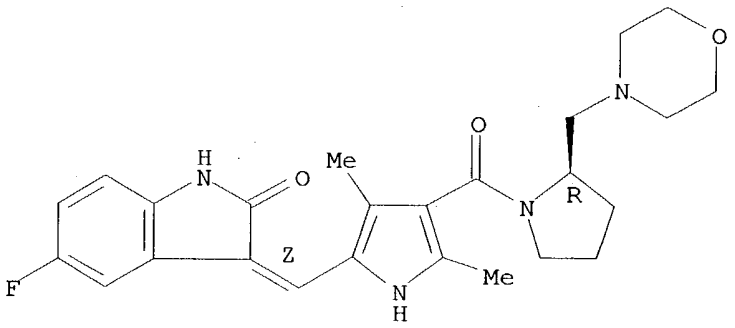
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Pyrrolidine, 1-[[5-[(Z)-(5-fluoro-1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-2-(4-morpholinylmethyl)-, (2R)- (9CI)
MF C25 H29 F N4 O3

Absolute stereochemistry.
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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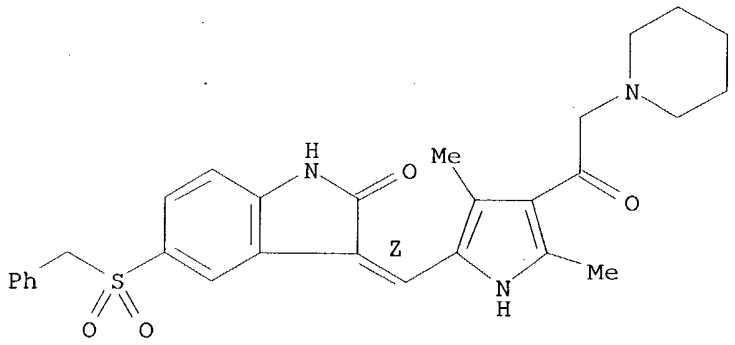
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 3-[[3,5-dimethyl-4-(1-piperidinylacetyl)-1H-pyrrol-2-yl]methylene]-1,3-dihydro-5-[(phenylmethyl)sulfonyl]-, (3Z)- (9CI)
MF C29 H31 N3 O4 S

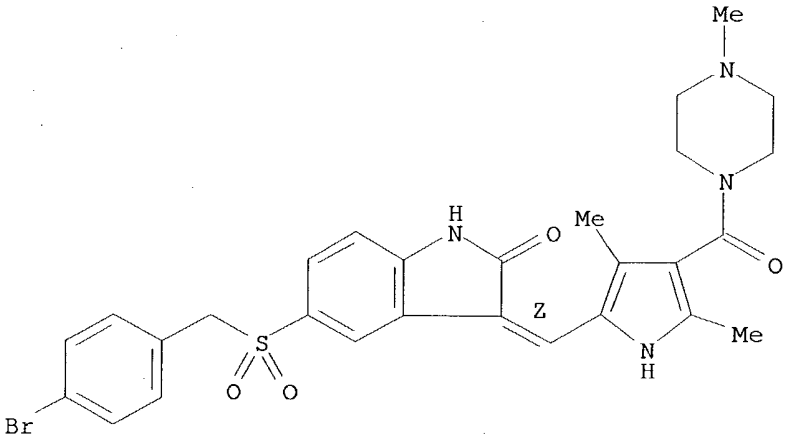
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperazine, 1-[[5-[(Z)-[5-[[(4-bromophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-4-methyl- (9CI)
MF C28 H29 Br N4 O4 S

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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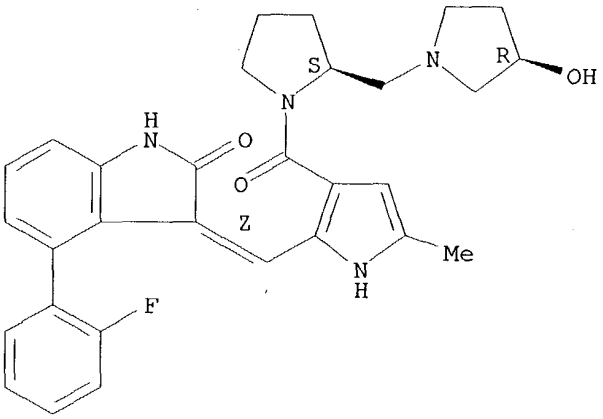
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Pyrrolidine, 1-[[2-[(Z)-[4-(2-fluorophenyl)-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-5-methyl-1H-pyrrol-3-yl]carbonyl]-2-[[(3R)-3-hydroxy-1-pyrrolidinyl]methyl]-, (2S)- (9CI)
MF C30 H31 F N4 O3

Absolute stereochemistry.
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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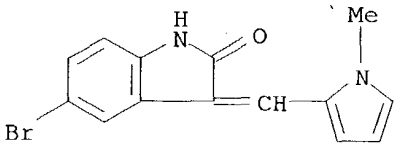
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 5-bromo-1,3-dihydro-3-[(1-methyl-1H-pyrrol-2-yl)methylene]-(9CI)
MF C14 H11 Br N2 O



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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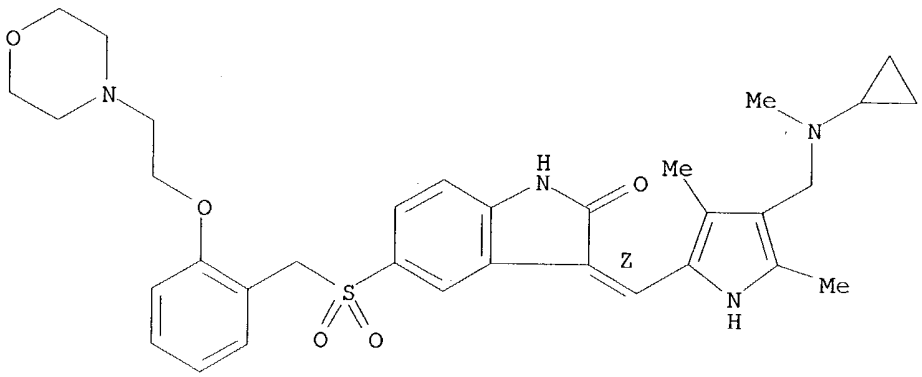
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 3-[[4-[(cyclopropylmethylamino)methyl]-3,5-dimethyl-1H-pyrrol-2-yl]methylene]-1,3-dihydro-5-[[[2-[2-(4-morpholinyl)ethoxy]phenyl]methyl]sulfonyl]-, (3Z)- (9CI)
MF C33 H40 N4 O5 S

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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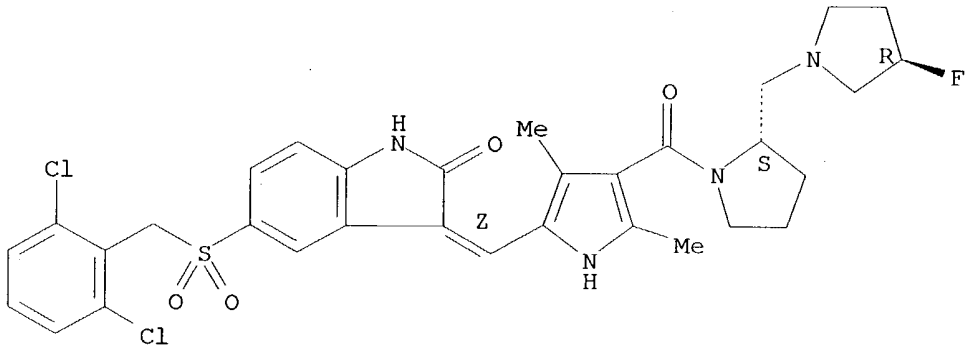
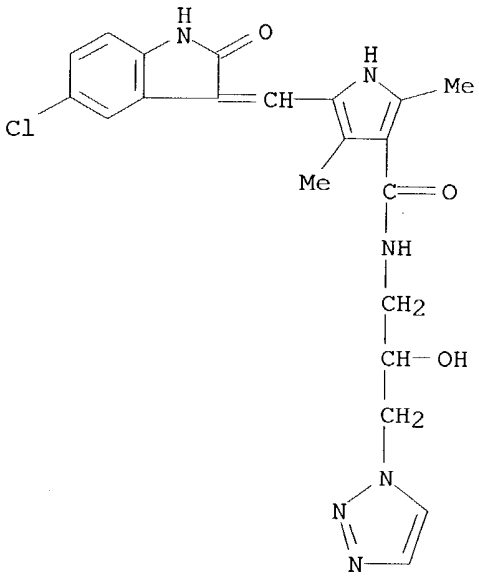
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Pyrrolidine, 1-[[5-[(Z)-[5-[[(2,6-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-2-[[(3R)-3-fluoro-1-pyrrolidinyl]methyl]-, (2S)- (9CI)
MF C32 H33 Cl2 F N4 O4 S

Absolute stereochemistry.
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-carboxamide, 5-[(5-chloro-1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-N-[2-hydroxy-3-(1H-1,2,3-triazol-1-yl)propyl]-2,4-dimethyl-(9CI)
MF C21 H21 Cl N6 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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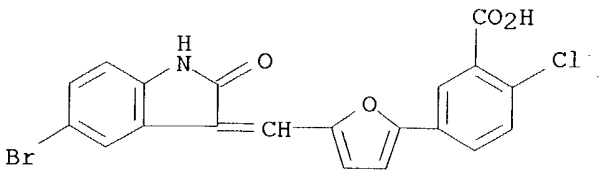
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Benzoic acid, 5-[5-[(5-bromo-1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-2-furanyl]-2-chloro- (9CI)
MF C20 H11 Br Cl N O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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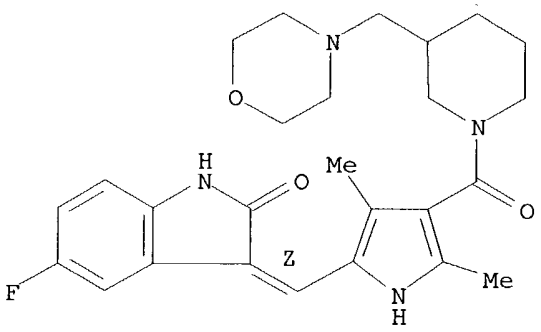
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperidine, 1-[[5-[(Z)-(5-fluoro-1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-3-(4-morpholinylmethyl)- (9CI)
MF C26 H31 F N4 O3

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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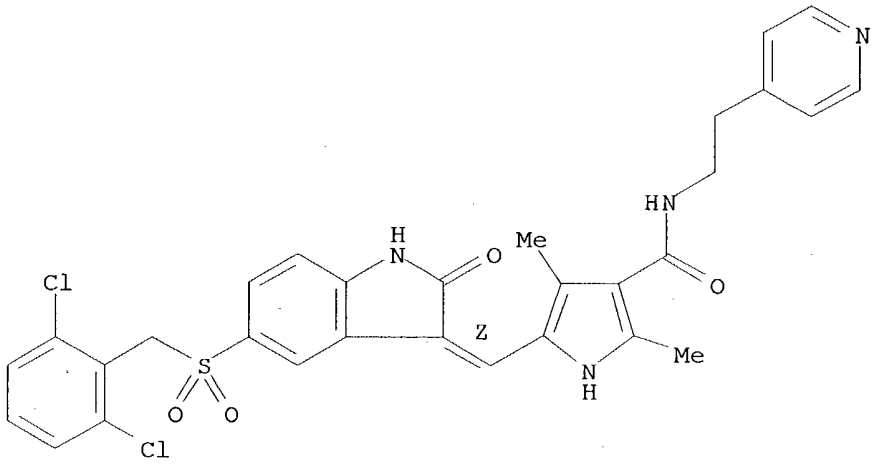
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-carboxamide, 5-[(Z)-[5-[[(2,6-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-N-[2-(4-pyridinyl)ethyl]- (9CI)
MF C30 H26 Cl2 N4 O4 S

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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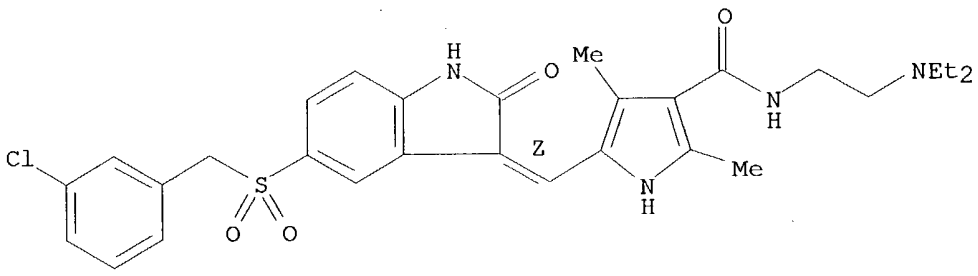
L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-carboxamide, 5-[(Z)-[5-[[(3-chlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-N-[2-(diethylamino)ethyl]-2,4-dimethyl- (9CI)
MF C29 H33 Cl N4 O4 S

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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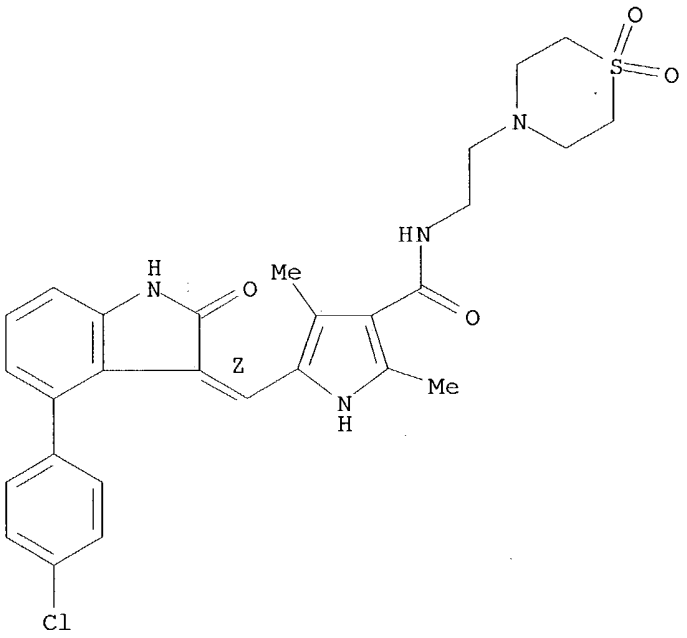
L3 50 ANSWERS - REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-carboxamide, 5-[(Z)-[4-(4-chlorophenyl)-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-N-[2-(1,1-dioxido-4-thiomorpholinyl)ethyl]-2,4-dimethyl- (9CI)
MF C28 H29 Cl N4 O4 S

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

V. Balasubramanian

=> log y

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.84	1.05

STN INTERNATIONAL LOGOFF AT 14:10:28 ON 24 AUG 2004